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**EP 0 404 190 B1**

CHEMICAL ABSTRACTS, vol. 103, no. 7, August 19, 1985, Columbus, Ohio, USA SIGOVA, V.I. et al. "Synthesis and biological activity of the aryl amides of 2-methylnicotinic and 2-phenylindolizine-8-carboxylic acids." page 563, column 1, abstract-no. 53 910y & Khim. Pharm. Zh. 1985, 19(3), 159-63

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## Description

## BACKGROUND OF THE INVENTION

## 5 1. Field of the Invention

This invention relates to novel condensed heterocyclic compounds or salts thereof, intermediates thereof, method for preparing them and farming and gardening fungicide containing them.

## 10 2. Description of the Prior Art

It has been known that some kinds of amide compounds possess fungicidal activity against pathogenic fungi causing various kinds of plant blight. For instance, compounds described in Japanese Patent Unexamined Publication No. 135,364/1988, EP 61836A and GB 2190375A exhibit effects against downy mildew fungi of a vine and cucumber and pathogenic fungi of a tomato and potato. However, it is hard to say that those compounds exhibit sufficient controlling effects and safety.

Sulfonylurea compounds which contain a condensed heterocyclic group and their herbicidal use are disclosed in EP 238 070.

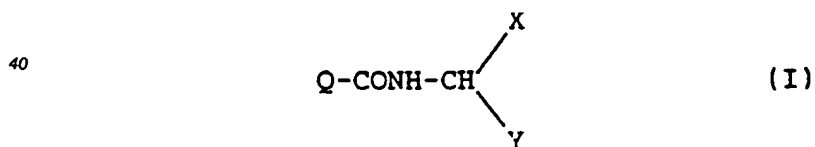
Thus, it is desired to conquer the aforementioned defects, namely, to provide compounds possessing excellent control effects against plant blight, method for preparing those compounds and agent for controlling plant blight containing those compounds.

Downy mildew and other blight of vegetables and fruit trees are liable to happen and spread in a time of much rain. Accordingly, it is required to provide compounds which are not washed away by rain (rain-resisting quality) and exhibit excellent effect for controlling and curing plant blight, do not damage the cultivating plants and show less toxicity to warm-blooded animals and fish, and also to provide a simple and easy method for preparing those compounds in high yield and to provide a useful agricultural chemical containing those compounds.

## SUMMARY OF THE INVENTION

The present inventors have studied intensively to attain the aforementioned problems, and succeeded in preparation of condensed heterocyclic compounds of the following general formula (I) and their salts and found that these compounds exhibit excellent effects for controlling various kinds of plant blight, especially downy mildew and blight, possess excellent rain-resisting quality, do not damage the plants and are less toxicity against warm-blooded animals and fish and that those compounds can be prepared by an industrially advantageous method.

Thus, the present invention relates to a condensed heterocyclic compound of the general formula:

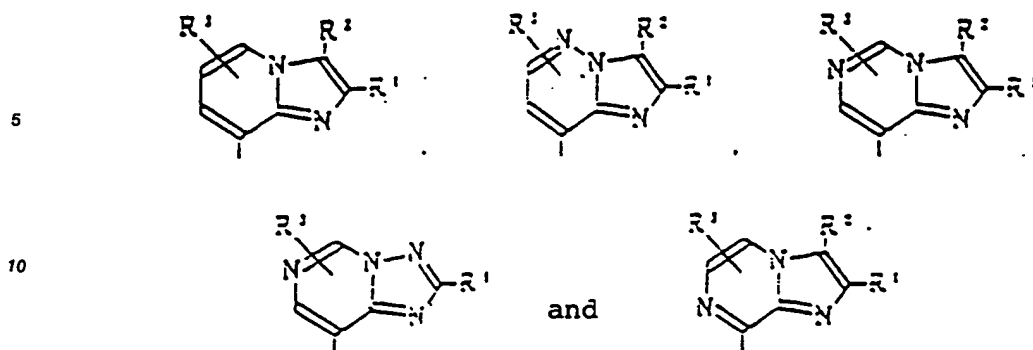


45 or an agriculturally acceptable salt thereof, wherein

(1) Q is a condensed heterocyclic group having a nitrogen atom in the bridgehead selected from the group consisting of

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15 in which R<sup>1</sup> is C<sub>1-6</sub> alkyl, halogen, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio, C<sub>6-10</sub> aryloxy, C<sub>6-10</sub> arylthio, alkoxycarbonyl, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group;

20 R<sup>2</sup> and R<sup>3</sup> are hydrogen, C<sub>1-6</sub> alkyl, halogen, nitro, amino, sulfo, mono- or dialkylsulfamoyl, alkoxycarbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group;

25 X is hydrogen; a group attached through a carbon atom selected from the group consisting of C<sub>1-10</sub> alkyl, C<sub>2-4</sub> alkenyl, C<sub>1-10</sub> haloalkyl, C<sub>3-6</sub> cycloalkyl, C<sub>3-6</sub> cycloalkenyl, C<sub>6-10</sub> aryl, C<sub>7-10</sub> aralkyl and a heterocyclic ring of the formula



35 a group attached through an oxygen atom selected from the group consisting of C<sub>1-10</sub> alkoxy, C<sub>6-10</sub> aryloxy and C<sub>7-10</sub> aralkyloxy; a group attached through a sulfur atom selected from the group consisting of C<sub>1-10</sub> alkylthio, C<sub>6-10</sub> arylthio and C<sub>7-10</sub> aralkylthio; a group attached through a nitrogen atom selected from the group consisting of C<sub>1-10</sub> alkylamino, C<sub>6-10</sub> arylamino, C<sub>7-10</sub> aralkylamino and a heterocyclic ring of the formula

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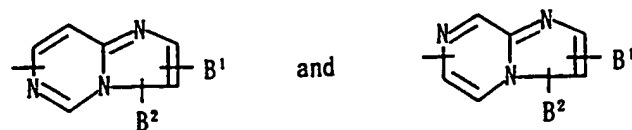
50 wherein each of said groups attached through a carbon atom, an oxygen atom, a sulfur atom or a nitrogen atom is unsubstituted or substituted with one to four substituents selected from the group consisting of nitro, amino, hydroxy, cyano, carboxyl, sulfonyl, halogen, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio, phenylthio and benzylthio; and Y is cyano, carbamoyl, thiocarbamoyl or trichloromethyl group;

55 or (2) Q is a condensed heterocyclic group having a nitrogen atom in the bridgehead selected from the group consisting of

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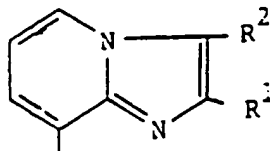


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in which the substituents B¹ and B² on the condensed heterocyclic group for Q may be the same or different each other, and are hydrogen atom, nitro, amino, hydroxyl, cyano, C₁-₃ acyl, carbamoyl, carboxyl, C₁-₄ alkoxy, sulfo, halogen, C₁-₄ alkoxy, C₆-₁₀ aryloxy, C₆-₁₀ arylcarbonyl, C₆-₁₀ aryl, C₇-₁₀ aralkyl, C₃-₇ cycloalkyl, C₁-₄ alkylthio, C₆-₁₀ arylthio, C₇-₁₀ aralkylthio, mono- or di-C₁-₄ alkylsulfamoyl, C₁-₆ alkyl, mono-, di-, tri- or tetra-halogen-substituted phenyl, mono-, di-, tri- or tetra-C₁-₄ alkyl-substituted phenyl, mono-, di-, tri- or tetra-C₁-₄ alkoxy-substituted phenyl, mono-, di-, tri- or tetra-C₁-₄ alkylthio-substituted phenyl, 2-chloro-4-nitrophenyl, 4-nitrophenyl, 2-methyl-4-aminophenyl, 2-bromo-4-nitrophenyl, 2-nitro-4-methylphenyl, or 5 or 6 membered heterocyclic group selected from pyridyl, furyl, thienyl and thiazolyl; and X and Y have the same meaning as defined above; or (3) Q is:

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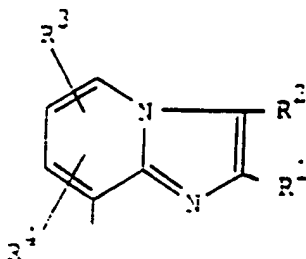
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and

R¹ is methyl, R² is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 or R¹ is methyl, R² is 1,1,2,2-tetrafluoroethyl, X is 2-furyl and Y is cyano;  
 or R¹ is methyl, R² is N,N-dimethylaminomethyl, X is 2-thienyl and Y is cyano;  
 or R¹ is methyl, R² is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R¹ is methyl, R² is morpholinomethyl, X is 2-furyl and Y is cyano;  
 or R¹ is methyl, R² is isopropylthiomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R¹ is methyl, R² is 2-cyanoethyl, X is 3-fluorophenyl and Y is cyano;  
 or R¹ is methyl, R² is 2-methyl-1-propenyl, X is 3-fluorophenyl and Y is cyano;  
 or R¹ is phenyl, R² is hydrogen, X is 2-N-methylpyrrolyl and Y is cyano;  
 or R¹ is phenyl, R² is hydrogen, X is 3-trifluoromethylphenyl and Y is cyano;  
 or R¹ is phenyl, R² is hydrogen, X is 3-methylphenyl and Y is cyano;  
 or R¹ is phenyl, R² is hydrogen, X is 5-methylfuryl-2-yl and Y is cyano;  
 or R¹ is phenyl, R² is hydrogen, X is 2-pyridylmethyl and Y is cyano;  
 or R¹ is phenyl, R² is hydrogen, X is 3-methylthien-2-yl and Y is cyano;  
 or R¹ is phenyl, R² is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 or R¹ is phenyl, R² is N,N-dimethylaminomethyl, X is 2-thienyl and Y is cyano;  
 or R¹ is phenyl, R² is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R¹ is phenyl, R² is 1,1,2,2-tetrafluoroethyl, X is 2-furyl and Y is cyano;  
 or R¹ is phenyl, R² is 2-cyanoethyl, X is 2-furyl and Y is cyano;  
 or R¹ is phenyl, R² is 2-methyl-1-propenyl, X is 2-furyl and Y is cyano;  
 or R¹ is phenyl, R² is 2,2,2-trichloro-1-hydroxyethyl, X is 2-furyl and Y is cyano;  
 or R¹ is trifluoromethyl, R² is hydrogen, X is 2-furyl and Y is cyano;  
 or R¹ is 2-methyl-1-propenyl, R² is hydrogen, X is 2-thienyl and Y is cyano;  
 or R¹ is ethoxyiminomethyl, R² is hydrogen, X is 2-thienyl and Y is cyano;  
 or R¹ is phenoxyiminomethyl, R² is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R¹ is phenylthiomethyl, R² is hydrogen, X is 3-fluorophenyl and Y is cyano;

55

or R<sup>1</sup> is phenylsulfinylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenylsulfonylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenoxyethyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is methoxymethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 5 or R<sup>1</sup> is methoxymethyl, R<sup>2</sup> is morpholinomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylthiomethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylsulfinylmethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylthiomethyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylsulfonylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 10 or R<sup>1</sup> is morpholinomethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is diethoxyphosphoxymethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is (dimethylthiocarbamoyl)thiomethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is chloromethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is trifluoromethyl, R<sup>2</sup> is phenylcarbamoyl, X is 3-fluorophenyl and Y is cyano;  
 15 or R<sup>1</sup> is hydrogen, R<sup>2</sup> is 1-hydroxybenzyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is sulfomethyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 4-dimethylaminophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is benzylidenamino, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is biphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 20 or R<sup>1</sup> is 4-sulfophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is trifluoromethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is formamide;  
 or R<sup>1</sup> is 3-hydroxyphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-hydroxyphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is benzylidenamino, X is 2-thienyl and Y is cyano;  
 25 or R<sup>1</sup> is 2-dimethylcarbamoylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methoxymethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-4-methylsulfinylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-4-methylsulfonylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 4-cyanophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 30 or R<sup>1</sup> is 2-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 3-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-5-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or (4) Q is:



and

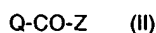
R<sup>1</sup> is methyl, R<sup>2</sup> is N,N-dimethylaminomethyl, R<sup>3</sup> is 5-methyl, R<sup>4</sup> is hydrogen, X is 2-thienyl and Y is cyano;

or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, R<sup>3</sup> is 5-methyl, R<sup>4</sup> is 7-methyl, X is 2-furyl and Y is cyano;

or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, R<sup>3</sup> is 5-trifluoromethyl, R<sup>4</sup> is 7-trifluoromethyl, X is 2-furyl and Y is cyano;

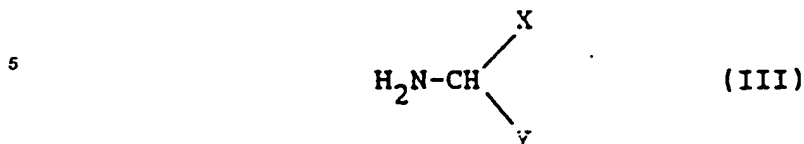
or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, R<sup>3</sup> is 5-hydroxy, R<sup>4</sup> is 6-chloro, X is 2-thienyl and Y is cyano.

This invention also relates to a method for preparing a condensed heterocyclic compound (I) or its salt by reacting a compound of the general formula:



wherein Q has the same meaning as defined in the above formula (I), and COZ denotes a carboxylic group

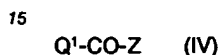
or its reactive derivative, or its salt with a compound of the general formula:



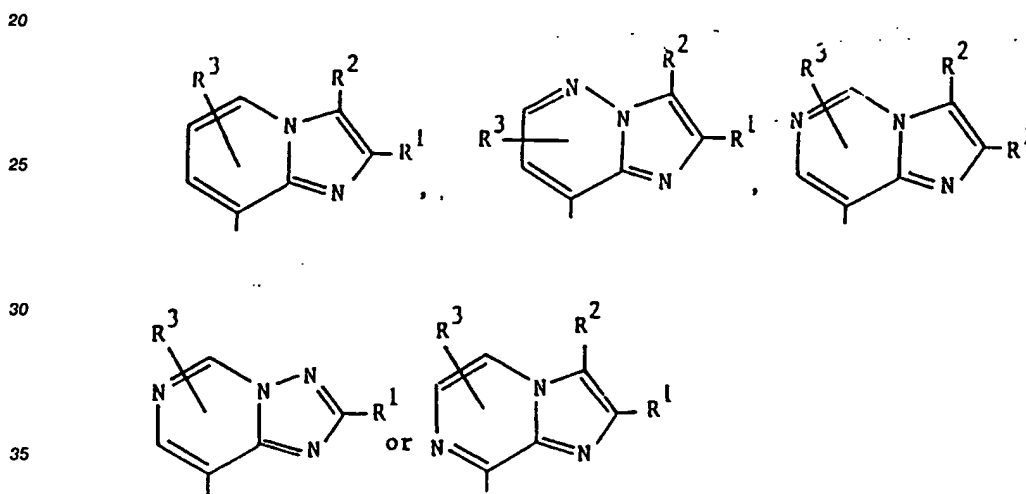
10 wherein X and Y are each as defined in the above formula (I), or its salt.

And further, this invention relates to a farming and gardening fungicide containing a condensed heterocyclic compound of the general formula (I) or its salt.

According to one aspect, the present invention provides a condensed heterocyclic carboxylic acid compound of the general formula:

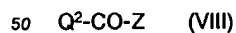


wherein Q<sup>1</sup> is a condensed heterocyclic group having a nitrogen atom in the bridgehead as follows



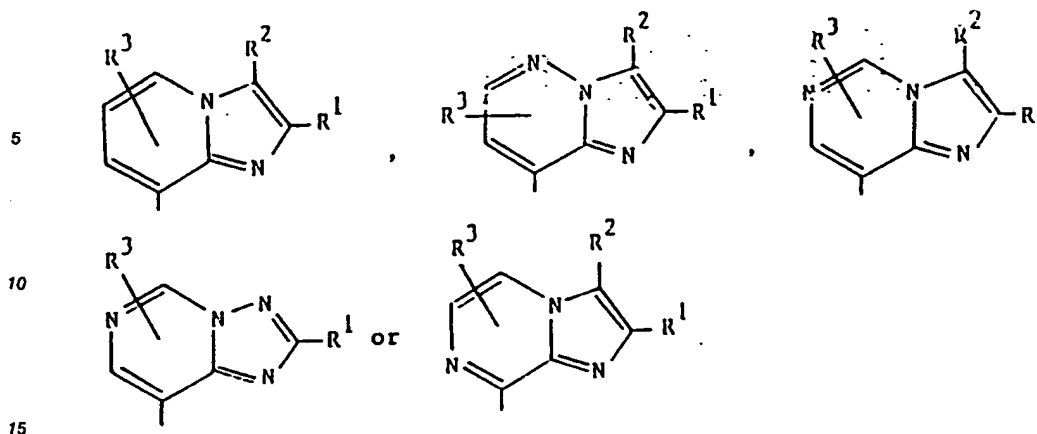
40 in which R<sup>1</sup> is an C<sub>1-6</sub> alkyl, halogen, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio, C<sub>6-10</sub> aryloxy, C<sub>6-10</sub> arylthio, alkoxycarbonyl, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group; R<sup>2</sup> and R<sup>3</sup> are a hydrogen, C<sub>1-6</sub> alkyl, halogen, nitro, amino, sulfo, mono- or dialkylsulfamoyl, alkoxycarbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group, and COZ denotes a carboxylic group or its reactive derivative; or its salt.

45 According to another aspect, the present invention provides a process for preparing a compound of the general formula:

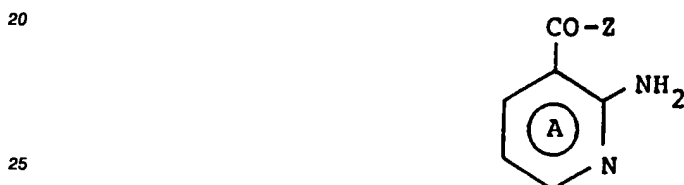


wherein Q<sup>2</sup> is a group of the general formula:

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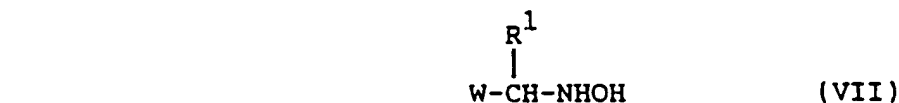
in which R<sup>1</sup>, R<sup>2</sup> and R<sup>3</sup> have the same meanings as defined above or its salt which comprises reacting a compound of the general formula:



wherein the ring  $\textcircled{A}$  is a pyridine, pyridazine, pyrimidine or pyrazine group which is unsubstituted or substituted by an C<sub>1-6</sub> alkyl, halogen, nitro, amino, sulfo, mono- or di-alkylsulfamoyl, alkoxycarbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group, and COZ denotes a carboxylic group or its reactive derivative, or its salt with a compound of the general formula:



40 or



45 wherein W is a halogen atom, and R<sup>1</sup> and R<sup>2</sup> have the same meanings as defined above.

The compounds (I) or their salts of this invention have a novel structure characterized by the combination of a specific group, namely, a condensed heterocyclic group having nitrogen atom in the bridgehead and the carbonyl group, which is different from the known amide compounds. The compounds of this invention have merits that they exhibit an excellent control effect for downy mildew and blight, possess excellent water-resisting quality, do not damage the plants and are less toxicity against warm-blooded animals and fish.

#### PREFERRED EMBODIMENT OF THE INVENTION

55 In the formula (I) and (II), the symbol Q has the above-described meaning. Examples of the substituent R<sup>1</sup> are alkoxycarbonyl (e.g. C<sub>1-4</sub> alkoxycarbonyl such as methoxycarbonyl, ethoxycarbonyl, n-propoxycarbonyl, isopropoxycarbonyl, etc.), halogen (chlorine, bromine, iodine, fluorine, etc.), C<sub>1-4</sub> alkoxy (methoxy,

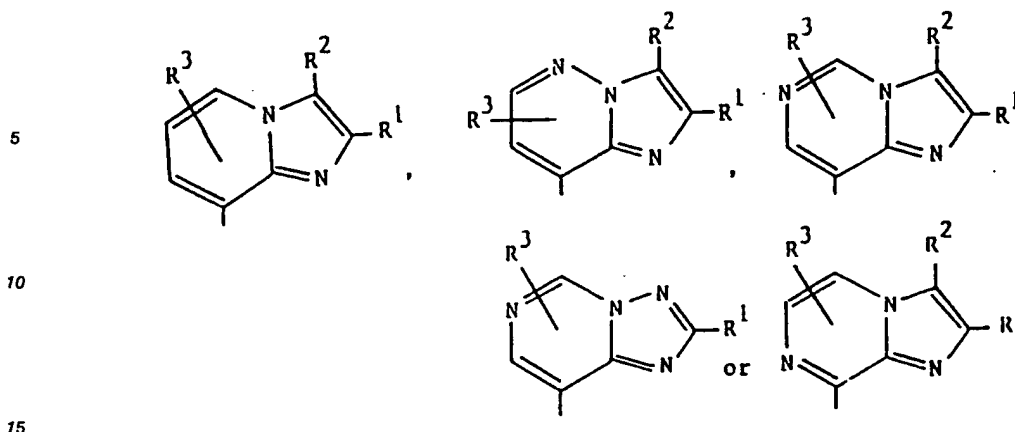


ethoxy, propoxy, isopropoxy, butoxy, isobutoxy, sec-butoxy, tert-butoxy, etc.), C<sub>6-10</sub> aryloxy (phenoxy, etc.), C<sub>1-4</sub> alkylthio (methylthio, ethylthio, propylthio, isopropylthio, butylthio, iso-butylthio, sec-butylthio, tert-butylthio, etc.), C<sub>6-10</sub> arylthio (phenylthio, etc.), C<sub>1-6</sub> alkyl (methyl, ethyl, propyl, isopropyl, pentyl, hexyl, etc.), phenyl, substituted phenyl (mono-, di-, tri or tetra-halogen-substituted phenyl such as 2-chlorophenyl, 3-chlorophenyl, 4-chlorophenyl, 2-fluorophenyl, 4-bromophenyl, 2,3-dichlorophenyl, 2,4-dichlorophenyl, 2,6-dichlorophenyl, 2,4-difluorophenyl, 2,6-difluorophenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkyl-substituted phenyl such as 2-methylphenyl, 3-methylphenyl, 4-ethylphenyl, 4-isopropylphenyl, 2,4-dimethylphenyl, 2,5-dimethylphenyl, 3,4-dimethylphenyl, 2,5-diethylphenyl, 2,4,6-trimethylphenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkoxy-substituted phenyl such as 2-methoxyphenyl, 3-methoxyphenyl, 4-methoxyphenyl, 3,4-dimethoxyphenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkylthio-substituted phenyl such as 2-methylthiophenyl, 3,4-dimethylthiophenyl, etc., 2-chloro-4-nitrophenyl, 4-nitrophenyl, 2-methyl-4-aminophenyl, 2-bromo-4-nitrophenyl, 2-nitro-4-methylphenyl, etc.), aromatic heterocyclic group (5 or 6 membered heterocyclic group such as pyridyl, furyl, thienyl, thiazolyl, etc.).

The substituents R<sup>2</sup> and R<sup>3</sup> on the condensed heterocyclic group for Q may be the same or different each other. Examples of the substituents are hydrogen atom, nitro, amino, cyano, formyl, alkoxycarbonyl (e.g., C<sub>1-4</sub> alkoxycarbonyl such as methoxycarbonyl, ethoxycarbonyl, n-propoxycarbonyl, isopropoxycarbonyl, etc.), sulfo, halogen (chlorine, bromine, iodine, fluorine, etc.), mono- or di-alkylsulfamoyl (e.g., mono- or di-C<sub>1-4</sub> alkylsulfamoyl such as mono- or di-methylsulfamoyl, mono- or di-ethylsulfamoyl, mono- or di-n-propylsulfamoyl), C<sub>1-6</sub> alkyl (methyl, ethyl, propyl, isopropyl, pentyl, hexyl, etc.), phenyl, substituted phenyl (mono-, di-, tri or tetra-halogen-substituted phenyl such as 2-chlorophenyl, 3-chlorophenyl, 4-chlorophenyl, 2-fluorophenyl, 4-bromophenyl, 2,3-dichlorophenyl, 2,4-dichlorophenyl, 2,6-dichlorophenyl, 2,4-difluorophenyl, 2,6-difluorophenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkyl-substituted phenyl such as 2-methylphenyl, 3-methylphenyl, 4-ethylphenyl, 4-isopropylphenyl, 2,4-dimethylphenyl, 2,5-dimethylphenyl, 3,4-dimethylphenyl, 2,5-diethylphenyl, 2,4,6-trimethylphenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkoxy-substituted phenyl such as 2-methoxyphenyl, 3-methoxyphenyl, 4-methoxyphenyl, 3,4-dimethoxyphenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkylthio-substituted phenyl such as 2-methylthiophenyl, 3,4-dimethylthiophenyl, etc., 2-chloro-4-nitrophenyl, 4-nitrophenyl, 2-methyl-4-aminophenyl, 2-bromo-4-nitrophenyl, 2-nitro-4-methylphenyl, etc.), aromatic heterocyclic group (5 or 6 membered heterocyclic group such as pyridyl, furyl, thienyl, thiazolyl, etc.).

The substituents B<sup>1</sup> and B<sup>2</sup> on the condensed heterocyclic group for Q may be the same or different each other. Examples of the substituents are hydrogen atom, nitro, amino, hydroxyl, cyano, C<sub>1-3</sub> acyl (formyl, etc.), carbamoyl, carboxyl, alkoxycarbonyl (e.g., C<sub>1-4</sub> alkoxycarbonyl such as methoxycarbonyl, ethoxycarbonyl, n-propoxycarbonyl, isopropoxycarbonyl, etc.), sulfo, halogen (chlorine, bromine, iodine, fluorine, etc.), C<sub>1-4</sub> alkoxy (methoxy, ethoxy, propoxy, isopropoxy, butoxy, isobutoxy, sec-butoxy, tert-butoxy, etc.), C<sub>6-10</sub> aryloxy (phenoxy, etc.), C<sub>6-10</sub> arylcarbonyl (benzoyl, etc.), C<sub>6-10</sub> aryl (phenyl, etc.), C<sub>7-10</sub> aralkyl (benzyl, phenethyl, etc.), C<sub>3-7</sub> cycloalkyl (cyclohexyl, etc.), C<sub>1-4</sub> alkylthio (methylthio, ethylthio, propylthio, isopropylthio, butylthio, iso-butylthio, sec-butylthio, tert-butylthio, etc.), C<sub>6-10</sub> arylthio (phenylthio, etc.), C<sub>7-10</sub> aralkylthio (benzylthio, etc.), mono- or di-alkylsulfamoyl (e.g., mono- or di-C<sub>1-4</sub> alkylsulfamoyl such as mono- or di-methylsulfamoyl, mono- or di-ethylsulfamoyl, mono- or di-n-propylsulfamoyl), C<sub>1-6</sub> alkyl (methyl, ethyl, propyl, isopropyl, pentyl, hexyl, etc.), substituted phenyl (mono-, di-, tri or tetra-halogen-substituted phenyl such as 2-chlorophenyl, 3-chlorophenyl, 4-chlorophenyl, 2-fluorophenyl, 4-bromophenyl, 2,3-dichlorophenyl, 2,4-dichlorophenyl, 2,6-dichlorophenyl, 2,4-difluorophenyl, 2,6-difluorophenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkyl-substituted phenyl such as 2-methylphenyl, 3-methylphenyl, 4-ethylphenyl, 4-isopropylphenyl, 2,4-dimethylphenyl, 2,5-dimethylphenyl, 3,4-dimethylphenyl, 2,5-diethylphenyl, 2,4,6-trimethylphenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkoxy-substituted phenyl such as 2-methoxyphenyl, 3-methoxyphenyl, 4-methoxyphenyl, 3,4-dimethoxyphenyl, etc., mono-, di-, tri- or tetra-C<sub>1-4</sub> alkylthio-substituted phenyl such as 2-methylthiophenyl, 3,4-dimethylthiophenyl, etc., 2-chloro-4-nitrophenyl, 4-nitrophenyl, 2-methyl-4-aminophenyl, 2-bromo-4-nitrophenyl, 2-nitro-4-methylphenyl, etc.), aromatic heterocyclic group (5 or 6 membered heterocyclic group such as pyridyl, furyl, thienyl, thiazolyl, etc.). Especially, one of B<sup>1</sup> and B<sup>2</sup> is preferably C<sub>1-6</sub> alkyl, phenyl or substituted phenyl.

Q<sup>1</sup> in the above mentioned general formula (IV) means a condensed heterocyclic group having a nitrogen atom in the bridgehead shown by the following formula:



The substituents  $R^1$ ,  $R^2$  and  $R^3$  have the same meanings as mentioned above.

Examples of the substituent  $R^1$  include the  $C_{1-6}$  alkyl, halogen,  $C_{1-4}$  alkoxy,  $C_{1-4}$  alkylthio,  $C_{6-10}$  aryloxy,  $C_{6-10}$  arylthio, alkoxycarbonyl, phenyl, substituted phenyl or aromatic heterocyclic group as stated above. Examples of the substituents  $R^2$  and  $R^3$  include hydrogen, nitro, amino, sulfo, formyl, cyano or phenyl group, and also the  $C_{1-6}$  alkyl, halogen, mono- or di-alkylsulfamoyl, alkoxycarbonyl, substituted phenyl or aromatic heterocyclic groups as stated above. The substituent  $R^1$  is preferably phenyl or a substituted phenyl.

Preferable ones of  $Q^1$  are  $Q'$  or  $Q''$  as described below.

25 The symbol X in the general formula (I) is a hydrogen atom or a group attached through C, O, S or N atom. The group attached through a C atom may be an alkyl, alkenyl, haloalkyl, cycloalkyl, cycloalkenyl, aryl or aralkyl group or an aromatic heterocyclic group having a chemical bond at a carbon atom; the group attached through an O atom may be an alkoxy, aryloxy or aralkyloxy group; the group attached through a S atom may be an alkylthio, arylthio or aralkylthio group; and the group attached through a N atom may be an alkylamino, arylamino or aralkylamino group or an aromatic heterocyclic group having a chemical bond at a N atom.

35 The above alkyl group and alkyl moiety of the haloalkyl, alkoxy, alkylthio and alkylamino groups include a straight or branched one containing 1 to 10 carbon atoms such as methyl, ethyl, propyl, isopropyl, n-butyl, isobutyl, sec-butyl, tert-butyl, n-pentyl, sec-pentyl, isopentyl, neopentyl, n-hexyl, isohexyl, n-octyl, n-decyl and the like.

The above alkenyl group may be one containing 2 to 4 carbon atoms such as vinyl, allyl, 2-methylallyl, 3-methylallyl, 3-butenyl or the like.

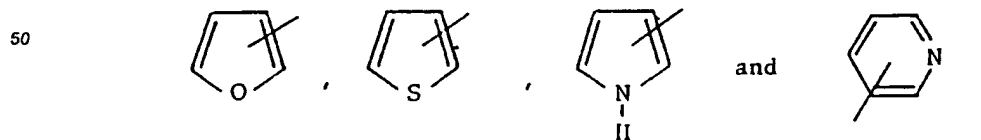
The above cycloalkyl group may be a 3- to 6-membered one such as cyclopropyl, cyclopentyl, cyclohexyl or the like.

40 The above cycloalkenyl group may be a 3- to 6-membered one such as cyclopropenyl, cyclopentenyl, cyclohexenyl or the like.

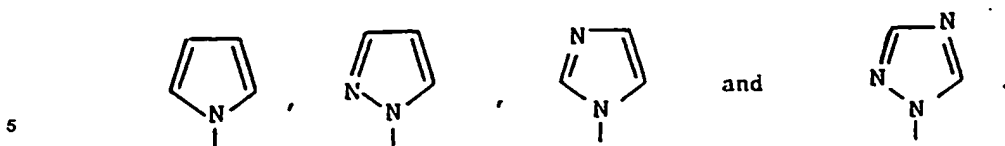
The above aryl group and aryl moiety of the aryloxy, arylthio and arylamino groups may be one containing 6 to 10 carbon atoms such as phenyl, tolyl, xylyl, naphthyl or the like.

45 The above aralkyl group and aralkyl moiety of the aralkyloxy, aralkylthio and aralkylamino groups may be one containing 7 - 10 carbon atoms such as benzyl, phenethyl or the like.

The aromatic heterocyclic group having a chemical bond at a C atom may be a 5 or 6-membered aromatic heterocyclic one such as



The aromatic heterocyclic group having a chemical bond at a N atom may be a 5 or 6-membered N-containing one such as

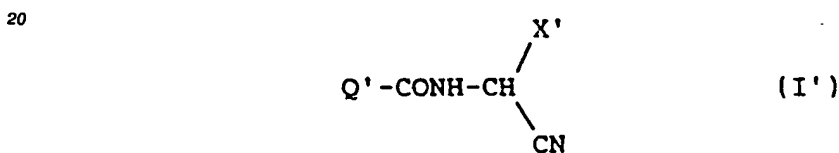


Each of the above mentioned groups attached through C, O, S or N atom may further have one to four substituents such as nitro, amino, hydroxyl, cyano, carboxyl, sulfo, a halogen (fluorine, chlorine, bromine, etc.), an alkoxy containing 1 to 4 carbon atoms (methoxy, ethoxy, etc.), an alkylthio containing 1 to 4 carbon atom(s) (methylthio, ethylthio, etc.), phenylthio, benzylthio and the like.

Suitable examples of X include hydrogen atom, an  $C_{1-10}$  alkyl,  $C_{2-4}$  alkenyl,  $C_{1-10}$  alkoxy,  $C_{1-10}$  alkylthio, phenyl, halogen-substituted phenyl, aromatic heterocyclic group having a chemical bond at C or N atom or the like. Specific examples of these groups include the ones exemplified for  $R^1$ ,  $R^2$  and  $R^3$  in the above.

The symbol Y in the general formula (I) means an electron attractive group, such as cyano, carbamoyl, thiocarbamoyl or trichloromethyl group or the like. Cyano group is preferable as an example of Y.

Especially interesting compounds of this invention are ones represented by the following formula (I'):

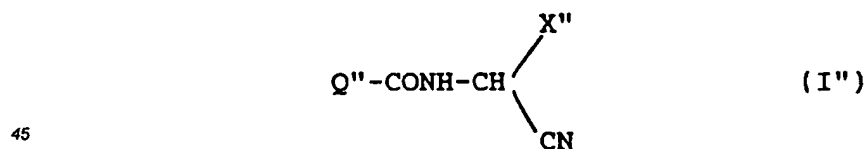


wherein  $Q'$  is a group of the formula:



35 (one of  $B^1$  and  $B^2$  is an  $C_{1-6}$  alkyl, phenyl or substituted phenyl group and the other is a hydrogen atom),  $X'$  is a phenyl, halogen-substituted phenyl or aromatic heterocyclic group having a chemical bond at a C atom.

Another group of interesting compounds (I) is one represented by the following formula (I''):



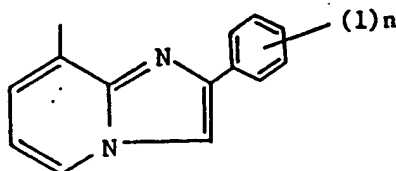
wherein  $Q''$  is



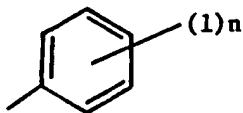
(B<sup>1''</sup> is a phenyl or substituted phenyl group, B<sup>2''</sup> is a hydrogen atom or a C<sub>1-6</sub> alkyl group), and X' is a phenyl, fluorine-substituted phenyl, thienyl or furyl group.

The C<sub>1-6</sub> alkyl group and substituted phenyl group exemplified in the above R<sup>1</sup>, R<sup>2</sup> and R<sup>3</sup> are applicable to those in B<sup>1'</sup>, B<sup>1''</sup>, B<sup>2'</sup> and B<sup>2''</sup> of the formulas (I') and (I''). The halogen-substituted phenyl group for X' may be a phenyl group substituted by one to four halogens of fluorine, chlorine, bromine and so on. The fluorine-substituted phenyl group for X'' may be o-, m- or p-fluorophenyl, o,m- or o,p-difluorophenyl or 2,4,6-trifluorophenyl. The aromatic heterocyclic groups having a chemical bond at C atom mentioned in the symbol X are applicable to those of X'.

Preferable example of the symbols Q' and Q'' is



wherein n is 0, 1, 2, 3 or 4 and I is the same or different selected from a group comprising C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio or halogen. The substituted phenyl groups mentioned in R<sup>1</sup>, R<sup>2</sup> and R<sup>3</sup> are applicable to the group:



(n 0). X' and X'' is preferably 2-thienyl group, etc.

The compounds(I) of this invention contain an asymmetric carbon atom, and this invention include each isomer due to the asymmetric carbon atom and a mixture thereof.

The salt of the compounds (I) of this invention may be a salt formed by an alkali metal or alkaline earth metal such as sodium, magnesium, potassium, calcium, etc. with an acidic group (e.g., carboxy, etc.) contained as a substituent, or a salt formed by a mineral acid such as hydrochloric acid, phosphoric acid, sulfuric acid, etc. or an organic acid such as oxalic acid, acetic acid, benzoic acid, etc. with a basic group contained in the substituent(s) or condensed heterocycle.

The carboxylic acids (II) and (IV) having condensed heterocycle, which are used as the starting materials in this invention, or their salts can be prepared by a method similar to the known methods described in, for example, J. Org. Chem., Vol. 37, page 3107 (1972); J. Org. Chem., Vol. 36, page 2678 (1971); J. Org. Chem., Vol. 42, page 4197 (1977); J. Med. Chem., Vol. 17, page 645 (1974); J. Med. Chem., Vol. 20, page 386 (1977); J. Med. Chem., Vol. 15, page 982 (1972); J. Med. Chem., Vol. 28, page 876 (1985); Tetrahedron Lett., Vol. 21, page 2195 (1980); J. Chem., Soc., Perkin Trans. I, page 1159 (1987), and the like.

In more detail, some of the carboxylic acids (II) or (IV) or their salts can be prepared, for example, by reacting an aromatic heterocyclic aminocarboxylic acid, its reactive derivative or salt with a compound of the formula (VI) or (VII).

Especially, a compound (VIII) or its salt which is frequently used among novel compounds (IV) or their salts can be prepared by reacting a compound (V) or its salt with a compound (VI) or a compound (VII).

In the formulas (II), (IV), (V) and (VIII), Z denotes a leaving group and accordingly -COZ can denote -COOH or its reactive derivative. Thus, the compound (II) or its salt can be represented by a carboxylic acid of the formula:

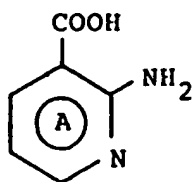


wherein Q has the same meaning as defined above, or its reactive derivative, or its salt; the compound (IV) or its salt be represented by a carboxylic acid of the formula:



wherein Q' has the same meaning as defined above, or its reactive derivative, or its salt;  
the compound (V) or its salt be represented by a carboxylic acid of the formula:

5



(V')

10

wherein the ring (A) has the same meaning as defined above  
or its reactive derivative, or its salt;

15 the compound (VIII) or its salt be represented by a carboxylic acid of the formula:



wherein  $Q^2$  has the same meaning as defined above, or its reactive derivative, or its salt, respectively.

20 Examples of the reactive derivatives at the carboxyl group in the aromatic heterocyclic amino carboxylic acid and carboxylic acids (II'), (IV'), (V') and (VIII') include the acid anhydrides, active amides, active esters or the like. Specific examples of such reactive derivatives are as follows:

1) Acid anhydrides may be mixed anhydrides with halogenic acids (e.g., hydrochloric acid or hydrobromic acid), mono-alkyl carbonates, aliphatic carboxylic acids (e.g., acetic acid, pivalic acid, 25 valeric acid, isovaleric acid or trichloroacetic acid) or aromatic carboxylic acids (e.g., benzoic acid), or symmetric acid anhydrides.

2) Active amides may be amides with pyrazole, imidazole, 4-substituted imidazole, dimethylpyrazole or benzotriazole.

3) Active esters may be methyl ester, ethyl ester, methoxymethyl ester, propargyl ester, 4-nitrophenyl ester, 2,4-dinitrophenyl ester, trichlorophenyl ester, pentachlorophenyl ester or mesitylphenyl ester, or 30 other esters with 1-hydroxy-1H-2-pyrrolidone, N-hydroxysuccinimide or N-hydroxyphthalimide.

Such reactive derivatives at carboxyl group are suitably selected and used depending upon the kind of the used carboxylic acids.

Further, the above mentioned aromatic heterocyclic amino carboxylic acid and carboxylic acids (II'), (IV'), (V') and (VIII') may be used in their free form (i.e., carboxylic acid form), preferably in the presence of 35 a condensing agent such as N,N'-dicyclohexylcarbodiimide, N-cyclohexyl-N'-morpholinoethylcarbodiimide, N-cyclohexyl-N'-(4-dimethylaminocyclohexyl)carbodiimide or N-ethyl-N'-(3-dimethylaminopropyl)-carbodiimide. In addition, the above carboxylic acids may be used as the salts with an alkali metal or alkaline earth metal.

40 The symbol W in the formula (VI) and (VII) denotes a halogen atom such as chlorine, bromine or fluorine. The ring (A) in the formula (V) may be pyridine, pyridazine, pyrimidine or pyrazine ring, which may be substituted by a group as mentioned in R<sup>3</sup>,  $Q^2$  is preferably Q' or Q''.

This reaction can be carried out without solvent or in a suitable solvent and in the presence of a 45 condensing agent such as a base, if necessary. The suitable solvents may be, alcohols such as ethanol, isopropyl alcohol, etc., aromatic hydrocarbons such as benzene, toluene, xylene, etc., halogenated hydrocarbons such as methylene chloride, chloroform, carbon tetrachloride, etc., ethers such as diethyl ether, dioxane, tetrahydrofuran, dimethoxyethane, etc., ketones such as acetone, methyl ethyl ketone, etc., nitriles such as acetonitrile, propionitrile, etc., amides such as dimethylformamide, dimethylacetamide, etc., esters 50 such as methyl acetate, ethyl acetate, butyl acetate, etc., or a mixture thereof. If necessary, a mixed solvent of water and an aromatic hydrocarbon or halogenated hydrocarbon may be used. The solvent is usually used in 1 to 50 times (weight), preferably 5 to 10 times, to the carboxylic acids or their reactive derivatives or salts. The reaction may be accelerated by adding a base to the solvent. Suitable base may be, tertiary amines such as triethylamine, pyridine, 4-dimethylaminopyridine, DBU (1,8-diazabicyclo[5.4.0]undec-7- 55 ene), etc., alkali metal or alkaline earth metal hydroxide, carbonate or bicarbonate or alkali metal alcoholate such as sodium carbonate, potassium carbonate, sodium bicarbonate, sodium hydroxide, potassium hydroxide, sodium alcoholate, etc., or organic metal salt such as n-butyl lithium, etc. The base is usually used in 1 to 5 equivalents, preferably, 1 to 3 equivalents, to the carboxylic acids or their reactive derivatives

or salts.

The compound of the formula (VI) or (VII) is used in an amount of 1.1 to 2.0 moles, to 1 mole of the carboxylic acid or its reactive derivative or salt in this reaction. This reaction may be carried out at room temperature (e.g., -10 °C to 30 °C) or at an elevated temperature (e.g., 30 to 100 °C) for accelerating the reaction. The reaction time depends on the reaction temperature, etc., and it is usually 15 minutes to 24 hours, preferably 30 minutes to 10 hours.

Thus, the resulting compound (II) or (IV) can be isolated and purified as a free base by a known method such as chromatography on silica gel (Kiesel gel® 60 manufactured by Merck & Co., Inc., eluting solvent: chloroform, ethyl acetate, etc.), or isolated and purified as an acid addition salt with an inorganic acid such as hydrochloric acid, sulfuric acid, phosphoric acid, etc., or with an organic acid such as acetic acid, benzenesulfonic acid, p-toluenesulfonic acid, methanesulfonic acid, citric acid, tartaric acid, oxalic acid, propionic acid, maleic acid, malic acid, malonic acid, fumaric acid, mandelic acid, ascorbic acid, etc. according to a conventional method by conventional means such as concentration, concentration under reduced pressure, extraction, phase transfer, crystallization, recrystallization or chromatography. In case where a carboxy group or the like is contained as a substituent in the compound, said compound can be derived to a salt with alkali metal or alkaline earth metal as exemplified before according to a conventional method, which can be further isolated and purified by conventional means as mentioned above.

The condensed heterocyclic carboxylic acid derivative (II) or (IV) or its salt prepared by the above reaction can also be derived to a nitro-substituted compound by a conventional nitration, or to a halogenated compound by using a halogenating agent such as N-bromosuccinimide (NBS), N-chlorosuccinimide (NCS), or the like.

The amines (III) and their salts (salts with the acids as mentioned in the above compounds (II) and (IV)) can be prepared by a known method or method similar thereto.

A compound (I) or its salt can be prepared by reacting a compound (II) or its salt with a compound (III) or its salt.

In more detail, a compound (I) or its salt can usually be prepared by reacting a compound (II) or its salt with a compound (III) or its salt without solvent or in a suitable solvent and in the presence of a suitable base or condensing agent. The suitable solvents may be, aromatic hydrocarbons such as benzene, toluene, xylene, etc., halogenated hydrocarbons such as methylene chloride, chloroform, carbon tetrachloride, etc., ethers such as diethyl ether, dioxane, tetrahydrofuran, etc., ketones such as acetone, methyl ethyl ketone, etc., nitriles such as acetonitrile, propionitrile, etc., amides such as dimethylformamide, dimethylacetamide, etc., esters such as methyl acetate, ethyl acetate, butyl acetate, etc., or a mixture thereof. If necessary, a mixture of water and an aromatic hydrocarbon or halogenated hydrocarbon may be used. The solvent is usually used in 1 to 50 times (weight), preferably 5 to 10 times, to the compound (II) or its salt. The reaction can be accelerated by adding a base to the solvent. The suitable bases may be, tertiary amines such as triethylamine, pyridine, 4-dimethylaminopyridine, DBU (1,8-diazabicyclo[5.4.0]undec-7-ene), etc., alkali metal or alkaline earth metal hydroxides, carbonates or bicarbonates or alkali metal alcoholates such as sodium carbonate, potassium carbonate, sodium bicarbonate, sodium hydroxide, potassium hydroxide, sodium alcoholate etc., or organic metal salt such as n-butyl lithium, etc. Such base is usually used in 1 to 20 equivalents, preferably 2 to 5 equivalents, to the compound (II) or its salt. The suitable condensing agents may be, thionyl chloride, phosphorus oxy-chloride, carbonyldiimidazole, N-methyl-2-bromo-pyridinium iodide, dicyclohexylimide, etc., or a mixture thereof. Such condensing agent is usually used in 1 to 10 equivalents, preferably 2 to 5 equivalents, to the compound (II) or its salt.

In the reaction, the compound (II) or its salt is used in an amount of 1.1 to 1.5 moles to 1 mole of the compound (III) or its salt. The reaction may be carried out under cooling to at room temperature (-20 to 30 °C), or at an elevated temperature such as about 30 to 100 °C to accelerate the reaction. Though the reaction time depends the reaction temperature, etc., it is usually 15 minutes to 15 hours, preferably about 30 minutes to 8 hours.

Thus the resulting compound (I) can be isolated and purified as a free base by a known method such as chromatography on silica gel (Kiesel gel 60 manufactured by Merck & Co, Inc., eluting solvent: chloroform, ethyl acetate etc.), or isolated and purified as an acid addition salt with an inorganic acid such as hydrochloric acid, sulfuric acid, phosphoric acid, etc., or with an organic acid such as acetic acid, benzenesulfonic acid, p-toluenesulfonic acid, methanesulfonic acid, citric acid, tartaric acid, oxalic acid, propionic acid, maleic acid, malic acid, malonic acid, fumaric acid, mandelic acid, ascorbic acid, etc. according to a conventional method, by conventional means as mentioned above. In case where a carboxy group or the like is contained as a substituent in the compound, said compound can be derived to a salt with alkali metal or alkaline earth metal as exemplified before according to a conventional method, and can be isolated and purified by conventional means as mentioned above.

Thus prepared compounds (I) of this invention or their salts possess an excellent effect for preventing and controlling various kinds of plant diseases caused by pathogenic fungi, especially downy mildew of vegetables such as cucumber, Chinese cabbage, onion, pulse, etc. and fruit trees such as vine, citrus, apple, etc., plant blight of tomato, potato, egg plant, green pepper, pumpkin, etc. Further, the compounds (I) of this invention and their salts maintain the stable fungicidal effect for a considerably long period of time after applying to plants (long-lasting effect), and are less in reduction of the efficacy due to little washing off by rain after application by spraying (rain-resistance effect). Thus, the compounds (I) and their salts exert sufficient effects even in the rainy season when downy mildew or plant blight occurs frequently. And further, the compounds (I) of this invention and their salts possess a safe and advantageous qualities as an agricultural fungicide, because damage on plants is low and toxicity to fish is also low.

The compound (I) of this invention or its salt can be used as a fungicide in a conventional form of agricultural chemicals. That is, a kind or two or more kinds of the compounds (I) of this invention and their salts may preferably be used in accordance with purpose in a form of emulsifiable concentrate, oily preparation, spray, wettable powder, powder, tablet, ointment, etc., which can be formed with a suitable carrier or carriers in a conventional manner, for example, by dissolved or dispersed in a proper liquid carrier, by mixed with a proper solid carrier or adsorbed on a proper solid carrier. If necessary, emulsifying agent, suspending agent, spreading agent, penetrating agent, moisturizing agent, adhesive, stabilizing agent, etc. may be added to the aforementioned preparations, which can be prepared by a conventional method.

The suitable rate of the compound (I) of this invention or its salt to the total fungicidal composition is about 1 to 80% by weight for emulsifiable concentrate or wettable powder, about 0.1 to 10% by weight for oily preparation and powder, and about 5 to 50% by weight for granules, though the concentration of the active ingredient may be varied in accordance with the purpose. The emulsifiable concentrates, wettable powders or the like is suitably diluted and extended (for example to 100 - 5000 times) with water or the like on the occasion of use, and then scattered.

Suitable liquid carrier (solvent) to be used may be, for example, water, alcohols (e.g. methanol, ethanol, n-propanol, isopropyl alcohol, ethylene glycol, etc.), ethers (e.g. dioxane, tetrahydrofuran, ethylene glycol monomethyl ether, diethylene glycol monomethyl ether, propylene glycol monomethyl ether, etc.), ketones (e.g. acetone, methyl ethyl ketone, etc.), aliphatic hydrocarbons (e.g. kerosene, kerosene, fuel oil, machine oil, etc.), aromatic hydrocarbons (e.g. benzene, toluene, xylene, solvent naphtha, methylnaphthalene, etc.), halogenated hydrocarbons (e.g. methylene chloride, chloroform, carbon tetrachloride, etc.), acid amides (e.g. dimethylformamide, dimethylacetamide, etc.), esters (e.g. ethyl acetate, butyl acetate, fatty acid glycerol ester, etc.), nitriles (e.g. acetonitrile, propionitrile, etc.), or the like. These liquid carriers may be used individually or as an optional mixture thereof.

Suitable solid carrier (diluent or extender) may be vegetable powders (e.g. soybean flour, tobacco flour, wheat flour, sawdust, etc.), mineral powders (e.g. clays such as kaolin, bentonite, acid clay, talcs such as talc powder, pyrophyllite, silicas such as diatomaceous earth, mica powder, etc.), alumina, sulfur powder, active carbon, or the like. These solid carriers may be used individually or as an optional mixture thereof.

Suitable base for the ointment may be, for example, polyethylene glycol, pectin, polyalcohol esters of higher fatty acids such as glycerin mono-stearate, etc. cellulose derivatives such as methyl cellulose, etc., sodium alginate, bentonite, higher alcohols, polyalcohols such as glycerin, etc., vaseline, white petrolatum, liquid paraffin, lard, various kinds of vegetable oils, lanolin, dehydrated lanolin, hard oil, resins, or the like. These bases may be used individually or as an optional mixture thereof, or together with the surface active agent exemplified in the following.

Suitable surface active agent to be used, if necessary, as the emulsifying agent, spreading agent, penetrating agent, dispersing agent, etc. may be nonionic or anionic surface active agent such as soaps, polyoxyalkyl aryl esters (e.g. Nonal®, manufactured by Takemoto Yushi K.K., Japan), alkyl sulfates (e.g. Emal 10®, Emal 40®, manufactured by Kao-Atlas K.K., Japan), alkyl sulfonates (e.g. , Neogen®, Neogen T®, manufactured by Dai-ichi Kogyo Seiyaku K.K., Japan : Neopellex®, manufactured by Kao-Atlas K.K., Japan), polyethylene glycol ethers (e.g. Nonipol 85®, Nonipol 100®, Nonipol 160®, manufactured by Sanyo Kasei K.K. Japan), polyalcohol esters (e.g. Tween 20®, Tween 80®, manufactured by Kao-Atlas K.K., Japan), or the like.

The fungicide of this invention can be applied at any time from seeds nursery plant to harvest. The fungicide of this invention can be used for controlling the outbreak of plant blight by previous administration and also for curing the plant blight by administration after the outbreak according to a conventional method.

The amount of the compound (I) of this invention or its salt to be used as an agricultural chemicals is variable in accordance with conditions such as growth stage and growth situation of the plant to be applied, kind of the plant blight, situation of the outbreak, administration time and method of the fungicide, etc. The

compound (I) or its salt is usually used in an amount of about 3 to 300 g, preferably about 10 to 100 g per 10 are. The concentration of the active ingredient to be applied is preferably about 10 to 1000 ppm. The fungicide is usually applied by direct scattering, direct powdering or irrigating to the plant or by powdering to seeds. The amount, concentration and method of application may optionally be varied so far as the fungicide is used safely and effectively. The fungicide of this invention may be used, if necessary, in an optional combination with the other kind of fungicides (e.g. organic chlorine fungicide, organic phosphorus fungicide, benzimidazole-type fungicide, copper fungicide, organic sulfur fungicide, phenolic fungicide, triazole-type fungicide, pyrimidine-type fungicide, acrylic acid-type fungicide, sulfenamide-type fungicide, amino acid-type fungicide, antibiotic substance, etc.), insecticides (e.g. natural insecticide, carbamate-type insecticide, organic phosphorus insecticide, neaistoxin-type insecticide, synthesized pyrethroid, etc.), miticides, nematicides, herbicides, plant hormone drugs, plant growth regulators, stabilizing agents, synergists, attracting agents, repellents, perfumes, coloring agents, fertilizers, plant nutrients, various kinds of amino acids, low or high molecular phosphoric acids, or the like. Metal salts (e.g. copper chloride, copper sulfate, etc.) may also be added for the purpose of synergy.

The compounds (I) provided by this invention and their salts exhibit an excellent control or prevent effect against downy mildew and blight of vegetables and fruit trees. The compounds (I) or their salts are not washed away, possess an excellent rain-resistant property and accordingly exhibit the excellent control or prevent effect especially in a rainy season. The compounds (I) of this invention and their salts decrease the blight of vegetables and fruit trees, and can be used as an advantageous fungicide with no substantial damage to the crops.

#### Test 1 Prevention of blight of tomato

A compound of this invention was dissolved in dimethylformamide (final concentration : 1% by weight), xylene (final concentration : 0.02% by weight) and Tween 20® (final concentration 0.02% by weight) were added to the solution, and the mixture was diluted with water to a fixed concentration of the active ingredient. To this solution was added a spreading agent, Dyne® (manufactured by Takeda Chemical Industries, Ltd., containing 20% (w/w) of polyoxyethylene nonyl phenyl ether and 12% of calcium lignin sulfonate) in a rate of 0.05% (w/w) (final concentration) to give a scattering solution. The solution was sprayed to young seedlings (about 4 weeks seedling) of tomato to an extent that the solution fell in drops. After air-drying, a suspension of zoospores of pathogen of tomato blight (concentration : about  $10^5$ /ml) was inoculated by spraying. After the inoculation, the plants were kept in a humid room at 17°C for 5 days. The rate of area of the affected spots in the plant was surveyed, and the prevention value was shown in accordance with the following coefficients.

Prevention value 3 : affected area 0 - 5%  
 Prevention value 2 : affected area 6 - 15%  
 Prevention value 1 : affected area 16 - 30%  
 Prevention value 0 : affected area 31% or more.

The results are shown in Table 1.

#### Test 2 Prevention of downy mildew of cucumber.

A scattering solution containing a compound of this invention in a fixed concentration was prepared by the method described in the above Test 1, and the solution was sprayed to young seedlings (about 3 weeks seedling) of cucumber to an extent that the solution fell in drops. After air-drying, a suspension of zoospores of pathogen of downy mildew of cucumber (concentration : about  $10^5$ /ml) was inoculated by spraying. After the inoculation, the plants were kept in a humid room at 20°C for one day and further in a humid room for 6 days. The rate of area of the affected spots in the plant was surveyed, and the prevention value was shown in accordance with the following coefficients.

Prevention value 3 : affected area 0 - 5%  
 Prevention value 2 : affected area 6 - 15%  
 Prevention value 1 : affected area 16 - 30%  
 Prevention value 0 : affected area 31% or more

The results are shown in Table 2.



Test 3 Prevention of downy mildew of vine.

A scattering solution containing a compound of this invention in a fixed concentration was prepared by the method described in the above Test 1, and the solution was sprayed to young trees of vine (about 6 weeks seedling) to an extent that the solution fell in drops. After air-drying, a suspension of zoospores of pathogen of downy mildew of vine (concentration : about  $10^5$ /ml) was inoculated by spraying. After inoculation, the plants were kept in a humid room at 18°C for 10 days. The rate of area of the affected spots in the plant was surveyed, and the prevention value was shown in accordance with the following coefficients.

- 10    Prevention value 3 :    affected area 0 - 5%
- Prevention value 2 :    affected area 6 - 15%
- Prevention value 1 :    affected area 16 - 30%
- Prevention value 0 :    affected are 31% or more

The results are shown in Table 3.

- 15    In the following Tables 1 - 3, the numbers in the column of the test compound mean the compound No. obtained in the Examples described below.

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Table 1  
Prevention of  
blight of tomato

Compound No.	200ppm	Compound No.	200ppm
2-1	3	2-45	3
2-2	3		
2-3	3	2-52	3
2-4	3	2-54	3
2-5	3	2-55	3
2-7	3	2-56	3
2-9	3	2-57	3
2-10	3	2-58	3
2-11	2	2-59	3
2-15	3	2-63	3
2-18	3	2-64	3
2-19	3	2-65	3
2-20	3	2-66	3
2-24	3	2-67	3
2-25	3	2-68	3
2-26	3	2-69	3
2-27	3	2-70	2
2-32	3	2-71	3
2-33	3	2-72	3
2-34	3	2-73	3
2-35	2	2-78	3
2-41	2		

Table 1 (continued)

5	Compound No.	200ppm	Compound No.	200ppm
	2-79	3	2-145	3
	2-80	3	2-146	3
10	2-81	1	2-147	3
	2-82	3	2-148	3
	2-83	3	2-149	3
15	2-92	2	2-150	3
	2-104	3	2-151	3
	2-109	3	2-152	3
20	2-114	3	2-153	3
	2-115	2	2-154	3
	2-119	3	2-155	3
	2-120	3	2-156	3
25	2-121	3	2-157	3
	2-122	3	2-158	3
	2-123	3	2-159	3
30	2-127	3	2-160	3
	2-128	3	2-161	3
	2-134	3	2-162	3
35	2-135	3	2-163	3
	2-136	3	2-164	3
	2-137	3	2-166	3
	2-138	3	2-167	3
40	2-139	3	2-168	3
	2-140	3	2-170	3
	2-141	3	2-171	3
45	2-142	3	2-173	3
	2-143	3	2-174	3
	2-144	3	2-175	3
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Table 1 (continued)

Compound No.	200ppm	Compound No.	200ppm
2-176	3	2-190	3
2-177	3	2-194	3
2-178	3	2-195	3
2-179	3	2-196	3
2-180	3	2-197	3
2-181	3	2-198	3
2-182	3	2-200	3
2-183	3	2-201	3
2-184	3	2-203	3
2-185	3	2-204	3
2-186	3	2-211	3
2-187	3	2-212	3
2-188	3	2-214	3
2-189	3	2-215	3

Tabel 2  
Prevention of downy  
mildew of cucumber

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Compound No.	200ppm	Compound No.	200ppm
2-1	3	2-41	3
2-2	3	2-42	3
2-3	3	2-45	3
2-4	3		
2-5	3	2-48	3
2-7	3	2-52	3
2-9	3	2-54	3
2-10	3	2-55	3
2-11	3	2-56	3
2-15	3	2-57'	3
2-18	3	2-58	3
2-19	3	2-59	3
2-20	3	2-63	3
2-24	3	2-64	3
2-25	3	2-65	3
2-26	3	2-66	3
2-27	3	2-67	3
2-28	3	2-68	3
2-29	3	2-69	3
2-32	3	2-70	3
2-33	3	2-71	3
2-34	3	2-72	3
2-35	2	2-73	3
2-37	3	2-78	3
2-40	1	2-79	3

Table 2 (continued)

	Compound No.	200ppm	Compound No.	200ppm
5	2-81	3	2-148	3
	2-82	3	2-149	3
10	2-83	3	2-150	3
	2-92	3	2-151	3
	2-104	3	2-152	2
15	2-109	3	2-153	3
	2-114	3	2-154	3
	2-115	3	2-155	3
20	2-119	3	2-156	3
	2-120	3	2-157	3
	2-121	3	2-158	3
	2-122	3	2-159	3
25	2-123	3	2-160	3
	2-127	3	2-161	3
	2-128	3	2-162	3
30	2-134	3	2-163	3
	2-135	3	2-164	3
	2-136	3	2-166	3
35	2-137	3	2-167	3
	2-138	3	2-168	3
	2-139	3	2-170	3
40	2-140	3	2-171	3
	2-141	3	2-173	3
	2-142	3	2-174	3
	2-143	3	2-175	3
45	2-144	3	2-176	3
	2-145	3	2-177	3
	2-146	3	2-178	3
50	2-147	3	2-179	3

Table 2 (continued)

5	Compound No.	200ppm	Compound No.	200ppm
	2-180	3	2-195	3
	2-181	3	2-196	3
10	2-182	3	2-197	3
	2-183	3	2-200	3
	2-184	3	2-201	3
15	2-185	3	2-203	3
	2-186	3	2-204	3
	2-187	3	2-211	3
20	2-188	3	2-212	3
	2-189	3	2-214	3
	2-190	3	2-215	3
25	2-194	3		

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Table 3  
Prevention of downy  
mildew of vine

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Compound No.	200ppm	Compound No.	200ppm
2-1	3	2-35	3
2-2	3	2-36	3
2-4	3	2-37	3
2-5	3	2-41	3
2-7	3	2-42	1
2-9	2		
2-10	2	2-48	3
2-11	3	2-49	3
2-15	3		
2-18	2	2-54'	2
2-19	3	2-55	3
2-20	2	2-56	3
2-24	3	2-57	3
2-25	3	2-58	3
2-26	3	2-59	3
2-27	3	2-63	2
2-34	3	2-64	2



Table 3 (continued)

	Compound No.	200ppm	Compound No.	200ppm
5	2-65	3	2-143	3
	2-66	3	2-144	3
10	2-67	3	2-145	3
	2-68	2	2-146	3
	2-69	3	2-147	3
15	2-70	3	2-148	3
	2-71	2	2-149	3
	2-72	2	2-150	3
20	2-73	3	2-151	3
	2-78	2	2-152	3
	2-79	3	2-153	3
	2-82	3	2-154	3
25	2-83	3	2-155	3
	2-92	3	2-156	3
	2-104	1	2-157	3
30	2-109	3	2-158	3
	2-119	3	2-159	3
	2-120	3	2-160	3
35	2-121	3	2-161	3
	2-122	3	2-162	3
	2-123	3	2-163	3
	2-127	3	2-164	2
40	2-128	3	2-165	2
	2-134	2	2-166	3
	2-135	3	2-167	3
45	2-140	3	2-168	3
	2-141	3	2-170	3
	2-142	3	2-171	3
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It is shown by the above test results that the compounds (I) of this invention and their salts possess an excellent effect for preventing the downy mildew and blight of vegetables and fruit trees.

#### 55 (f) Examples

This invention is explained in more detail by the following Examples. However, this invention shall not be limited to these Examples.

In the following Examples, elution in the column chromatography was carried out under the observation of TLC (Thin Layer Chromatography). The observation of TLC was conducted by using Kesel gel 60F 254 (Art. 5715) manufactured by Merck & Co. as a TLC plate, the same solvent as the one used as an eluent of column chromatography as a developing solvent, and UV detector as a detecting method. The silica gel  
 5 filled in a column was Kesel gel 60 (70 - 230 mesh, Art. 7734) manufactured by Merck & Co. The NMR spectrum means proton NMR (<sup>1</sup>HNMR), and tetramethylsilane was used as an internal or external standard. The NMR spectrum was measured by VARIAN EM 390 (90 MHz) type spectrometer in case special description is not given, and each  $\delta$  value was shown by ppm. Numerical value in the parentheses means mixing rate by volume of each solvent used in a mixed solvent as an eluent.

10 Each abridgment used in the following Examples has the following meanings.

s : singlet, d : doublet, m : multiplet, br : broad, J : coupling constant, Hz : Hertz, CDCl<sub>3</sub> : heavy chloroform, DMSO-d<sub>6</sub> : heavy dimethyl sulfoxide, % (except for yield) : w/w %.

The room temperature means about 15 to 25 °C.

#### 15 Example 1

Synthesis of methyl imidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1 - 1)

47% Hydrogen bromide (2.5ml, 14.4m moles) was added to diethoxybromoethane (2.0ml, 13.2m  
 20 moles), and the mixture was stirred at 50 °C for 2 hours. The reaction mixture was cooled to room temperature, and ethanol (7.0ml) and sodium bicarbonate (1.0g, 11.9m moles) were added thereto. The mixture was stirred, and insoluble substance was filtered off. To the filtrate were added methyl 2-aminonicotinate (1.0g, 6.6m moles), sodium bicarbonate (2.0g, 13.8m moles) and ethanol (7.0ml), and the mixture was refluxed under heating for 4 hours. The reaction mixture was cooled to room temperature, and  
 25 saturated aqueous solution of sodium bicarbonate was added thereto. The mixture was extracted with dichloromethane (100ml x 3), and the extract was washed with water, dried over anhydrous magnesium sulfate (MgSO<sub>4</sub>) and concentrated under reduced pressure. The concentrate was purified by column chromatography on silica gel (eluent : chloroform and then chloroform/methanol=9/1) to give the object compound (0.6g, yield 55%) as yellow crystals. m.p. 67 - 69 °C.

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#### Example 2

Synthesis of methyl 2-methylimidazo[1,2-a]-pyridine-8-carboxylate (Compound No. 1 - 3)

35 Bromoacetone (6.5g, 42.7m moles) and methyl 2-aminonicotinate (4.8g, 31.5m moles) were added to ethanol (50ml), and the mixture was refluxed under heating for 17 hours. The reaction mixture was cooled to room temperature, and the solvent was distilled off under reduced pressure. Saturated aqueous solution of sodium bicarbonate (50ml) was added to the residue, and the mixture was extracted with chloroform (100ml x 3). The extract was washed with water, dried over MgSO<sub>4</sub> and concentrated under reduced pressure. The  
 40 concentrate was purified by column chromatography on silica gel (eluent:chloroform and then chloroform/methanol = 9/1) to give the object compound (4.0g, yield 85%) as an oily substance.

#### Example 3

45 Synthesis of hydrogen bromide salt of ethyl 2-phenyl-imidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1-25).

Phenacyl bromide (11.0g, 55.3m moles) and ethyl 2-aminonicotinate (8.68g, 52.2m moles) were added to methyl ethyl ketone (100 ml), and the mixture was refluxed under heating for 5.5 hours. The reaction  
 50 mixture was left to stand at room temperature overnight, and the precipitating crystals were collected by filtration to give the object compound (16.3g, yield 89.96%) as crystals. m.p. 174 - 176 °C.

#### Example 4

55 Synthesis of methyl 3-chloro-2-methylimidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1-15).

Methyl 2-methylimidazo[1,2-a]pyridine-8-carboxylate (5.3g, 27.9m moles) was dissolved in chloroform (30ml), and N-chlorosuccinimide (3.7g, 27.9m moles) was added to the solution and reacted at room

temperature for 40 minutes. To the reaction mixture was added 10% aqueous solution of sodium carbonate (80ml), and the mixture was stirred for 30 minutes. The organic layer was separated from the reaction mixture, and the aqueous layer was extracted with chloroform (10ml x 3). The extracts were combined, washed with water, dried over  $\text{MgSO}_4$  and concentrated under reduced pressure. The concentrate was purified by column chromatography on silica gel (eluent:chloroform) to give the object compound (3.6g, yield 57%) as an oily substance.

#### Example 5

10 Synthesis of methyl 3-dimethylaminomethyl-2-methylimidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1-17).

To acetonitrile (10ml) were added 37% formaldehyde (0.3ml) and acetic acid (0.4ml), and 50% aqueous solution of dimethylamine (0.3ml) was added slowly to the mixture at 0°C with stirring. Methyl 2-methylimidazo[1,2-a]pyridine-8-carboxylate (1.0g, 3.7m moles) was added to the mixture and reacted at 50°C for 2 hours and then at room temperature for 2.5 hours. The reaction mixture was concentrated under reduced pressure and neutralized with saturated aqueous solution of  $\text{NaHCO}_3$ . The solution was extracted with dichloromethane, and the extract was washed with water, dried over anhydrous sodium sulfate and concentrated under reduced pressure to give the object compound (0.48g, Yield 52%) as crystals. m.p. 123 - 124°C.

#### Example 6

Synthesis of 2-methylimidazo[1,2-a]pyridine-8-carboxylic acid (Compound No. 1-4)

25 Methyl 2-methylimidazo[1,2-a]pyridine-8-carboxylate (4g, 21m moles) was dissolved in a mixture of ethanol (45ml) and water (20ml), and sodium hydroxide (2.5g, 63m moles) was added thereto and refluxed under heating for 30 minutes. The reaction mixture was cooled to room temperature and concentrated under reduced pressure. Water (40ml) was added to the concentrate, and the mixture was adjusted to pH 4 with conc. hydrochloric acid. The precipitating crystals were collected by filtration to give the object compound (2.3g, yield 64%).

#### Example 7

35 Synthesis of 2-phenylimidazo[1,2-a]pyridine-8-carboxylic acid (Compound No. 1-26)

Hydrogen bromide salt of ethyl 2-phenylimidazo[1,2-a]pyridine-8-carboxylate (8.4g, 24.2m moles) was added to 20% aqueous solution of NaOH (200ml) and ethanol (100ml), and the mixture was refluxed under heating for 1 hour. The reaction mixture was cooled to room temperature, and alcohol was distilled off under reduced pressure. The concentrate was adjusted to pH 4 with conc. hydrochloric acid, and the precipitating crystals were collected by filtration to give the object compound (5.67g, yield 98.4%) as crystals. m.p. 226°C.

#### Example 8

45 Synthesis of 2-chloromethylimidazo[1,2-a]pyridine-8-carboxylic acid (Compound No. 1-90)

Ethyl 2-chloromethylimidazo[1,2-a]pyridine-8-carboxylate (8.4g, 35.2m moles) was added to conc. hydrochloric acid (100ml), and the mixture was refluxed under heating for 5 hours. The reaction mixture was cooled to room temperature and adjusted to pH 4 with saturated aqueous solution of sodium bicarbonate. The precipitating crystals were collected by filtration to give the object compound (5.67g, yield 76.9%) as crystal. m.p. 250°C (decomp.).

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## Example 9

## Synthesis of 3-nitro-2-(4-nitrophenyl)-imidazo[1,2-a]pyridine-8-carboxylic acid (Compound No. 1-96)

- 5 Ethyl 2-phenylimidazo[1,2-a]pyridine-8-carboxylate (2.9, 10m mol) was dissolved in conc. sulfuric acid (5ml), and to the solution was slowly added 70% nitric acid (1.3ml) at 10 °C. After addition, the mixture was stirred at room temperature for 30 minutes, poured into ice water and adjusted to pH 7 by adding 20% sodium hydroxide. The precipitated crystals were collected by filtration, which were ethyl ester of the title compound (3.6g, yield 99.2%). A solution of the crystals in conc. hydrochloric acid (35ml) was refluxed for 10 1.5 hours and cooled to room temperature. The precipitated crystals were collected by filtration and dried to give the object compound (2.9g, yield 87.3%). m.p. >300 °C.

## Example 10

- 15 Synthesis of 3-nitroimidazo[1,2-a]pyridine-8-carboxylic acid (Compound No. 1 - 125)

- Ethyl 2-ethoxycarbonylimidazo[1,2-a]pyridine-8-carboxylate (3.8g, 14.5m mol) was dissolved in conc. sulfuric acid (10ml), and to the solution was slowly added 70% nitric acid (1.5ml) at 10 °C. The mixture was stirred at room temperature for 30 minutes, poured into ice water and adjusted to pH 7 by adding 20% sodium hydroxide. The precipitated crystals were filtered to obtain ethyl 2-ethoxycarbonyl-3-nitro-imidazo- 20 [1,2-a]pyridine-8-carboxylate (2.4g, yield 53.8%) as crystals of m.p. 139 - 140 °C. A solution of the crystals in conc. hydrochloric acid (35ml) was refluxed for 1.5 hours and cooled to room temperature. The precipitated crystals were collected by filtration and dried to give the object compound (1.2g, yield 74.3%) of m.p. > 300 °C.

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## Example 11

## Synthesis of ethyl 3-formyl-2-phenyl-imidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1 - 119)

- 30 To a solution of ethyl 2-phenyl-imidazo[1,2-a]pyridine-8-carboxylate hydrogen bromide (14.7g, 4m mol) in DMF (8ml) was dropwise added phosphorus oxychloride (2ml) keeping the inner temperature to below 15 °C. Then, the mixture was allowed to react at room temperature for 30 minutes and at 70 °C for 2 hours. After completion of the reaction, the reaction mixture was cooled, poured into ice water and adjusted to pH 7 by adding 20% sodium hydroxide. The precipitated crystals were collected by filtration and recrystallized from ethanol to obtain the object compound (1.0g, yield 85.5%) as colorless crystals. m.p. 139 °C.

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## Example 12

## Synthesis of ethyl 3-cyano-2-phenyl-imidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1 - 121)

- 40 Ethyl 2-phenyl-imidazo[1,2-a]pyridine-8-carboxylate hydrogen bromide (1.96g, 5.35m mol), and triethylamine (0.55g, 5.44m mol) were added to acetonitrile (20ml). To this solution was dropwise added a solution of chlorosulfonylisocyanate (ClSO<sub>2</sub>NCO) (0.8ml, 9.19m mol) in acetonitrile (5ml) at an inner temperature of 0 - 20 °C. Then, the mixture was allowed to react at room temperature for an hour. After 45 completion of the reaction, the mixture was concentrated under reduced pressure. The resulting concentrate to which water was added was extracted with dichloromethane, and the extract was washed with water, dried over anhydrous sodium sulfate and concentrated under reduced pressure. The residue was purified by silica gel column chromatography (eluent : chloroform) to give the object compound (1.06g, yield 67.9%) as crystals. m.p. 87 - 89 °C.

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## Example 13

## Synthesis of ethyl 2-ethoxycarbonyl-imidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1 - 123)

- 55 A mixture of ethyl bromopyruvate (2.3g, 10.2m mol and ethyl 2-aminonicotinate (1.7g, 10m mol) in methyl ethyl ketone (17ml) was refluxed for 5 hours. After completion of the reaction, the mixture was cooled to room temperature and distilled under reduced pressure to remove the solvent. The residue to which saturated sodium hydrogen carbonate aqueous solution (50ml) was added was extracted with

chloroform (100ml x 3). The extract was washed with water, dried over  $\text{MgSO}_4$  and concentrated under reduced pressure. The concentrate was purified by silica gel column chromatography (eluent : ethyl acetate) to give the object compound (0.7g, yield 26.7%) as crystals. m.p.  $97^\circ\text{C}$ .

#### 5 Example 14

Synthesis of ethyl 2-methyl-3-sulfoimidazo[1,2-a]imidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1 - 127)

- 10 To a solution of ethyl 2-methyl-imidazo[1,2-a]pyridine-8-carboxylate (1.2g, 5.9m mol) in chloroform (15ml) was added chloresulfonic acid (0.6ml). The mixture was refluxed for 3 hours, and then concentrated under reduced pressure. The precipitated crystals were collected by filtration, washed sufficiently with water and dried to give the object compound (1.2g, yield 71.4%) as crystals. m.p.  $247 - 250^\circ\text{C}$  (decomp.).

#### 15 Example 15

Synthesis of ethyl 2-methyl-3-N-methylsulfamoylimidazo[1,2-a]pyridine-8-carboxylate (Compound No. 1 - 129)

- 20 Ethyl 2-methyl-3-sulfo-imidazo[1,2-a]pyridine-8-carboxylate (1.7g) and tri-n-propylamine (2.6g) were added to acetonitrile (17ml). Phosphorus oxychloride (1.1ml) was added to the mixture, keeping the inner temperature to  $50 - 60^\circ\text{C}$ . Then, the mixture was allowed to react at the same temperature for an hour. After completion of the reaction, the mixture was concentrated under reduced pressure, and water was added to the residue. The precipitated crystals were collected by filtration, washed with water and dried to obtain ethyl 2-methyl-3-chlorosulfonyl-imidazo[1,2-a]pyridine-8-carboxylate (1.2g, yield 66.3%) as crystals of m.p.  $182 - 183^\circ\text{C}$ .

- 25 The crystals (1.2g) was dissolved in acetonitrile (12ml), to which a solution of 40% methylamine aqueous solution (0.75g) in acetonitrile (3ml) was dropwise added at an inner temperature of  $0 - 10^\circ\text{C}$ . Then, the mixture was allowed to react at room temperature for an hour. After completion of the reaction, the mixture was concentrated under reduced pressure. The concentrate to which water was added was adjusted to pH 7 with dilute hydrochloric acid. The precipitated crystals was collected by filtration, washed with water and dried to obtain the object compound (1.1g, yield 93.2%) as crystals. m.p.  $184 - 185^\circ\text{C}$ .

#### Example 16

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Synthesis of 2-chloro-imidazo[1,2-a]pyridine-8-carboxylic acid (Compound No. 1 - 135)

- A mixture of diethyl bromomalonate (26ml) and ethyl 2-aminonicotinate (13.3g) was allowed to react at  $80 - 90^\circ\text{C}$  for 6 hours under nitrogen atmosphere. After completion of the reaction, the mixture was cooled to room temperature, and acetone (100ml) was added thereto. The precipitated crystals was collected by filtration, affording 3,8-diethoxycarbonyl-2-hydroxy-imidazo[1,2-a]pyridine hydrobromide (9.98g, yield 34.8%).

- 45 A mixture of the crystals (6.6g) in phosphorus oxychloride (30ml) was reacted at  $160^\circ\text{C}$  for 2 hours in a pressure reaction vessel. After completion of the reaction, excess phosphorus oxychloride was removed under reduced pressure. Ethanol was added to the residue, and the mixture was concentrated under reduced pressure. The concentrate was purified by silica gel column chromatography (eluent : chloroform) to obtain 2-chloro-3,8-diethoxycarbonyl-imidazo[1,2-a]pyridine (Compound No. 1 - 133) (2.2g, yield 40.3%) as crystals of m.p.  $105 - 106^\circ\text{C}$ .

- 50 The crystals (0.9g) were dissolved in ethanol (10ml), and 10% sodium hydroxide (5ml) was added thereto. The mixture was allowed to react at room temperature for an hour and then adjusted to pH 4 with conc. hydrochloric acid. The precipitated crystals were collected by filtration, washed with water and dried to obtain 2-chloro-imidazo[1,2-a]pyridine-3,8-dicarboxylic acid (0.73g, yield 100%) as crystals of m.p.  $210 - 212^\circ\text{C}$  (decomp.). Then, a mixture of the crystals (0.73g) in conc. hydrochloric acid (5ml) was refluxed for 2 hours. The reaction solution was adjusted to pH 4 with saturated sodium hydrogen carbonate aqueous solution, to yield crystals of the object compound (0.5g, yield 83.9%). m.p.  $237^\circ\text{C}$ .

## Example 17

Synthesis of  $\alpha$ -(imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-furyl)acetonitrile (Compound No. 2-1)

5  $\alpha$ -(2-Furyl)- $\alpha$ -aminoacetonitrile (0.7g, 5.5m mols) was dissolved in acetonitrile (20ml), and imidazo[1,2-a]pyridine-8-carboxylic acid chloride (1.0g, 5.5m moles) was added thereto with stirring and reacted at room temperature for 2 hours. After completion of the reaction, the solvent was removed from the reaction mixture under reduced pressure. To the residue were added water (20ml), saturated aqueous solution of sodium bicarbonate (50ml) and chloroform (50ml), and the mixture was stirred. The organic layer was  
 10 separated, dried over anhydrous magnesium sulfate and concentrated. The concentrate was purified by column chromatography on silica gel (eluent: ethyl acetate/n-hexane = 2/1) to give the object compound (0.16g, yield 11%) as crystals. m.p. 149 - 151 °C.

## Example 18

15 Synthesis of  $\alpha$ -(imidazo[1,2-a]pyridin-5-ylcarbonylamino)-(2-furyl)acetonitrile (Compound No. 2-224)

$\alpha$ -(2-Furyl)- $\alpha$ -aminoacetonitrile (0.7g, 5.5m moles) was dissolved in acetonitrile (20ml), and imidazo[1,2-a]pyridine-5-carboxylic acid chloride (1.0g, 5.5m moles) was added thereto and reacted at room temperature for 2 hours. After completion of the reaction, the solvent was removed from the reaction mixture. To the residue were added water (20ml), saturated aqueous solution of sodium bicarbonate (50ml) and chloroform (50ml), and the mixture was stirred. The organic layer was separated, dried over anhydrous magnesium sulfate and concentrated. The concentrate was purified by column chromatography on silica gel (eluent : methanol/chloroform = 1/9) to give the object compound (0.9g, yield 64%) as amorphous solid.

25  
 Example 19

Synthesis of  $\alpha$ -(2-phenyl-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-furyl)acetonitrile (Compound No. 2 - 24)

30 2-Phenyl-imidazo[1,2-a]pyridine-8-carboxylic acid (1.5g, 5.5m moles) was added to dried tetrahydrofuran (100ml), and carbonyldiimidazole (1.5g, 5.5m moles) was added slowly thereto with stirring at room temperature in an atmosphere of nitrogen gas. The mixture was stirred at room temperature for 8 hours, and  $\alpha$ -(2-furyl)- $\alpha$ -aminoacetonitrile (0.7g, 5.5m moles) was added to the mixture and reacted at room temperature for 5 hours. The solvent was removed from the reaction mixture under reduced pressure, and water (20ml) and ethyl acetate (50ml) were added to the residue and stirred. The organic layer was separated, dried over anhydrous magnesium sulfate and then concentrated. The concentrate was purified by column chromatography on silica gel (eluent : ethyl acetate/n-hexane = 1/1) to give the object compound (0.9g, yield 64%) as crystals. m.p. 197 - 199 °C.

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 Example 20

Synthesis of  $\alpha$ -(3-chloro-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-furyl)acetonitrile [Compound No. 2 - 10]

45 3-Chloro-imidazo[1,2-a]pyridine-8-carboxylic acid (1.8g, 9.2m moles) was added to thionyl chloride (20ml), and the mixture was refluxed under heating and stirring for 30 minutes and then concentrated under reduced pressure. The resultant crude acid chloride and  $\alpha$ -(2-furyl)- $\alpha$ -aminoacetonitrile (1.1g, 8.5m moles) were added to acetonitrile (50ml) and reacted at room temperature for 1 hour with stirring. The solvent was removed from the reaction mixture under reduced pressure, and water (20ml), saturated aqueous solution of sodium bicarbonate (50ml) and chloroform (50ml) were added to the residue and stirred. The organic layer was separated, dried over anhydrous magnesium sulfate and then concentrated. The concentrate was purified by column chromatography or silica gel (eluent : ethyl acetate/n-hexane = 1/2) to give the object compound (0.6g, yield 22%) as crystals. m.p. 142 - 143 °C.

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## Example 21

Synthesis of  $\alpha$ -(2-phenyl-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(3-chlorophenyl)acetonitrile (Compound No. 2 - 35)

5        2-Phenyl-imidazo[1,2-a]pyridine-8-carboxylic acid (1.0g, 4.2m moles),  $\alpha$ -(3-chlorophenyl)-aminoacetonitrile (0.77g, 4.6m moles) and  $\text{NaHCO}_3$  (1.1g, 13.1m moles) were added to acetonitrile (10ml). Phosphorus oxychloride (0.7ml, 7.51m moles, dissolved in 5ml of acetonitrile) was added slowly to the mixture with stirring under ice-cooling. The mixture was reacted at room temperature for 10 hours, and the  
10       solvent was removed from the reaction mixture under reduced pressure. Water (50ml) was added to the residue and the mixture was adjusted to pH 7 with saturated aqueous solution of sodium bicarbonate and then extracted with dichloromethane/ethyl acetate (10/1) (100ml x 2). The extract was dried over anhydrous magnesium sulfate and concentrated. Acetonitrile (5ml) was added to the concentrate, and the precipitating crystals were collected by filtration to give the object compound (0.44g, yield 27.2%) as crystals. m.p. 210 -  
15       211 °C.

## Example 22

20       Synthesis of  $\alpha$ -(3-dimethylaminomethyl-2-phenylimidazo[1,2-a]pyridin-8-ylcarbonylamino)-(3-fluorophenyl)-acetonitrile (Compound No. 2 - 79)

To acetonitrile (10ml) were added 37% formaldehyde (0.3ml) and acetic acid (0.4ml), and 50% aqueous solution of dimethylamine (0.3ml) was added slowly thereto at 0 °C with stirring. To the mixture was added  $\alpha$ -(2-phenyl-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(3-fluorophenyl)acetonitrile (1.0g, 2.7m moles), and the  
25       mixture was reacted at 50 °C for 2 hours and further stirred at room temperature for 2.5 hours. The reaction mixture was concentrated under reduced pressure, and the concentrate was neutralized with saturated aqueous solution of  $\text{NaHCO}_3$ . The solution was extracted with dichloromethane, and the extract was washed with water, dried over anhydrous sodium sulfate and concentrated under reduced pressure. The concentrate was purified by column chromatography or silica gel (eluent : chloroform) to give the object compound  
30       (0.47g, yield 40.8%) as crystals. m.p. 123 - 124 °C.

## Example 23

35       Synthesis of  $\alpha$ -(3-methylsulfinylmethyl-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-furyl)acetonitrile (Compound No. 2 - 122)

$\alpha$ -(2-Methylthiomethyl-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-furyl)acetonitrile (0.33g, 1m mole) was added to chloroform (10ml), and m-chloroperbenzoic acid (0.25g, 1.1m moles) was added to the mixture at internal temperature of 5 °C. The mixture was stirred at room temperature for 2.5 hours, and the  
40       reaction mixture was treated with saturated aqueous solution of  $\text{NaHCO}_3$ . The chloroform layer was separated, washed with water, dried over anhydrous sodium sulfate and then concentrated under reduced pressure. The concentrate was purified by column chromatography or silica gel (eluent : ethyl acetate) to give the object compound (0.1g, yield 30.3%) as crystals. m.p. 214 - 216 °C.

## 45       Example 24

Synthesis of  $\alpha$ -(2-(trifluoromethyl-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(3-fluorophenyl)acetamide (Compound No. 2 - 172)

50       A mixture of  $\alpha$ -(2-trifluoromethyl-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(3-fluorophenyl)acetonitrile (1.26g) and chlorosulfonic acid (0.5ml) in chloroform (13ml) was refluxed for 1.5 hours. The reaction solution was concentrated under reduced pressure, and water was added to the residue. The precipitated crystals were collected by filtration and recrystallized from ethanol to obtain the object compound (0.45g, yield 34.1%) as crystals. m.p. 118 - 120 °C

55       The compounds (II) obtained in a similar manner to those of Examples 1 - 16 are shown in Table 4, their melting points and NMR data are shown in Table 5, the compounds (I) obtained in a similar manner to those of Examples 17 - 24 are shown in Table 6, and their melting points and NMR data are shown in Table 7, respectively.

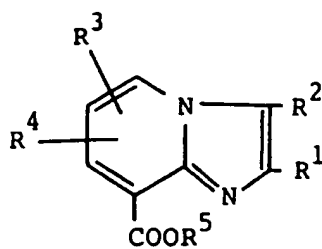




Table 4

	Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	1-1	H	H	H	H	Me
	1-2	H	H	H	H	H
	1-3	Me	H	H	H	Me
10	1-4	Me	H	H	H	H
	1-5	Et	H	H	H	Me
	1-6	Et	H	H	H	H
15	1-7	nPr	H	H	H	H
	1-8	iPr	H	H	H	Me
	1-9	iPr	H	H	H	H
20	1-10	nBu	H	H	H	H
	1-11	tBu	H	H	H	Me
	1-12	tBu	H	H	H	H
25	1-13	H	Cl	H	H	Et
	1-14	H	Cl	H	H	H
	1-15	Me	Cl	H	H	Me
30	1-16	Me	Cl	H	H	H
	1-17	Me	CH <sub>2</sub> NMe <sub>2</sub>	H	H	Me
	1-18	Me	CH <sub>2</sub> NMe <sub>2</sub>	H	H	H
35	1-19	Me	Me	H	H	H
	1-20	Me	CH <sub>2</sub> CH <sub>2</sub> CN	H	H	H
	1-21	Me	CHO	H	H	H
40	1-22	Me	NO <sub>2</sub>	H	H	H
	1-23	Me	CH=CHMe <sub>2</sub>	H	H	H
	1-24	C <sub>6</sub> H <sub>5</sub>	H	H	H	Me
45	1-25	C <sub>6</sub> H <sub>5</sub>	H	H	H	Et
	1-26	C <sub>6</sub> H <sub>5</sub>	H	H	H	H
	1-27	2-MeC <sub>6</sub> H <sub>4</sub>	H	H	H	Me

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Table 4 (continued)

Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	1-28	2-MeC <sub>6</sub> H <sub>4</sub>	H	H	H
	1-29	2-MeC <sub>6</sub> H <sub>4</sub>	H	H	Et
	1-30	3-MeC <sub>6</sub> H <sub>4</sub>	H	H	Me
10	1-31	3-MeC <sub>6</sub> H <sub>4</sub>	H	H	H
	1-32	4-MeC <sub>6</sub> H <sub>4</sub>	H	H	Me
	1-33	4-MeC <sub>6</sub> H <sub>4</sub>	H	H	H
15	1-34	2,5-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H
	1-35	2,4,6-Me <sub>3</sub> C <sub>6</sub> H <sub>2</sub>	H	H	H
	1-36	2-MeOC <sub>6</sub> H <sub>4</sub>	H	H	Me
20	1-37	2-MeOC <sub>6</sub> H <sub>4</sub>	H	H	H
	1-38	3-MeOC <sub>6</sub> H <sub>4</sub>	H	H	Me
	1-39	3-MeOC <sub>6</sub> H <sub>4</sub>	H	H	H
25	1-40	4-MeOC <sub>6</sub> H <sub>4</sub>	H	H	Me
	1-41	4-MeOC <sub>6</sub> H <sub>4</sub>	H	H	H
	1-42	3,4-(MeO) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	Et
30	1-43	3,4-(MeO) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H
	1-44	2-ClC <sub>6</sub> H <sub>4</sub>	H	H	Me
	1-45	2-ClC <sub>6</sub> H <sub>4</sub>	H	H	H
35	1-46	3-ClC <sub>6</sub> H <sub>4</sub>	H	H	Me
	1-47	3-ClC <sub>6</sub> H <sub>4</sub>	H	H	H
	1-48	4-ClC <sub>6</sub> H <sub>4</sub>	H	H	Me
40	1-49	4-ClC <sub>6</sub> H <sub>4</sub>	H	H	H
	1-50	4-BrC <sub>6</sub> H <sub>4</sub>	H	H	Me
	1-51	4-BrC <sub>6</sub> H <sub>4</sub>	H	H	H
45	1-52	4-NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub>	H	H	Me
	1-53	4-NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub>	H	H	H
	1-54	2-FC <sub>6</sub> H <sub>4</sub>	H	H	H

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Table 4 (continued)

	Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	1-55	2-F, 4-MeC <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-56	2-Me, 4-MeSC <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-57	2,4,6-Cl <sub>3</sub> C <sub>6</sub> H <sub>2</sub>	H	H	H	H
10	1-58	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> NMe <sub>2</sub>	H	H	Me
	1-59	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> NMe <sub>2</sub>	H	H	H
	1-60	C <sub>6</sub> H <sub>5</sub>	Cl	H	H	Me
15	1-61	C <sub>6</sub> H <sub>5</sub>	Cl	H	H	H
	1-62	C <sub>6</sub> H <sub>5</sub>	Me	H	H	H
	1-63	C <sub>6</sub> H <sub>5</sub>	C <sub>6</sub> H <sub>5</sub>	H	H	Me
20	1-64	C <sub>6</sub> H <sub>5</sub>	C <sub>6</sub> H <sub>5</sub>	H	H	H
	1-65	C <sub>6</sub> H <sub>5</sub>	CHO	H	H	Et
	1-66	C <sub>6</sub> H <sub>5</sub>	NO <sub>2</sub>	H	H	H
25	1-67	C <sub>6</sub> H <sub>5</sub>	CN	H	H	H
	1-68	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> CH <sub>2</sub> CN	H	H	H
	1-69	C <sub>6</sub> H <sub>5</sub>	CH=CMe <sub>2</sub>	H	H	H
30	1-70	C <sub>6</sub> H <sub>5</sub> CH <sub>2</sub>	H	H	H	H
	1-71	CF <sub>3</sub>	H	H	H	Et
	1-72	CF <sub>3</sub>	H	H	H	H
35	1-73	Cl	H	H	H	Et
	1-74	Cl	Cl	H	H	Et
	1-75	OMe	H	H	H	Et
40	1-76	OMe	H	H	H	H
	1-77	SMc	H	H	H	H
	1-78	H	OMe	H	H	H
45	1-79	2-thienyl	H	H	H	Et
	1-80	2-thienyl	H	H	H	H
50	1-81	C <sub>6</sub> H <sub>5</sub>	H	5-Me	H	Me

Table 4 (continued)

	Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	1-82	C <sub>6</sub> H <sub>5</sub>	H	5-Me	H	H
	1-83	C <sub>6</sub> H <sub>5</sub>	H	5-Me	7-Me	Et
	1-84	C <sub>6</sub> H <sub>5</sub>	H	5-Me	7-Me	H
10	1-85	C <sub>6</sub> H <sub>5</sub> OCH <sub>2</sub>	H	H	H	Me
	1-86	C <sub>6</sub> H <sub>5</sub> OCH <sub>2</sub>	H	H	H	H
	1-87	C <sub>6</sub> H <sub>5</sub> SCCH <sub>2</sub>	H	H	H	Et
15	1-88	C <sub>6</sub> H <sub>5</sub> SCCH <sub>2</sub>	H	H	H	H
	1-89	CH <sub>2</sub> Cl	H	H	H	Et
	1-90	CH <sub>2</sub> Cl	H	H	H	H
20	1-91	C <sub>6</sub> H <sub>5</sub> OCH <sub>2</sub>	CH <sub>2</sub> NMe <sub>2</sub>	H	H	Et
	1-92	CH <sub>2</sub> OMe	H	H	H	Et
	1-93	CH <sub>2</sub> SMc	H	H	H	Me
25	1-94	CH <sub>2</sub> SMc	H	H	H	H
	1-95	4-NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub>	NO <sub>2</sub>	H	H	Et
	1-96	4-NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub>	NO <sub>2</sub>	H	H	H
30	1-97	2,4-(NO <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	NO <sub>2</sub>	H	H	H
	1-98	2-thiazolyl	H	H	H	H
	1-99	4-Mc, 3-NO <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	NO <sub>2</sub>	H	H	H
35	1-100	2-Mc, 4-tBuC <sub>6</sub> H <sub>3</sub>	H	H	H	Et
	1-101	2-Mc, 4-tBuC <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-102	2-Mc, 5-tBuC <sub>6</sub> H <sub>3</sub>	H	H	H	Et
40	1-103	2-Mc, 5-tBuC <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-104	2,4-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-105	3,4-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	H
45	1-106	2,3,4,5-Me <sub>4</sub> C <sub>6</sub> H	H	H	H	Et
	1-107	2,3,4,5-Me <sub>4</sub> C <sub>6</sub> H	H	H	H	H

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Table 4 (continued)

	Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	1-108	2,5-Et <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	Et
	1-109	2,5-Et <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-110	4-ETC <sub>6</sub> H <sub>4</sub>	H	H	H	H
10	1-111	4-ETC <sub>6</sub> H <sub>4</sub>	H	H	H	Et
	1-112	2,5-Cl <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-113	2,5-(MeO) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	H
15	1-114	2-OMe,5-ClC <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-115	2-OMe,5-MeC <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-116	2-SMe,5-t-BuC <sub>6</sub> H <sub>3</sub>	H	H	H	H
20	1-117	2-SMe,5-MeC <sub>6</sub> H <sub>3</sub>	H	H	H	H
	1-118	2-Me,4-MeSC <sub>6</sub> H <sub>3</sub>	H	H	H	Et
	1-119	C <sub>6</sub> H <sub>5</sub>	CHO	H	H	Et
25	1-120	C <sub>6</sub> H <sub>5</sub>	CHO	H	H	H
	1-121	C <sub>6</sub> H <sub>5</sub>	CN	H	H	Et
	1-122	C <sub>6</sub> H <sub>5</sub>	CN	H	H	H
30	1-123	COOEt	H	H	H	Et
	1-124	COOEt	NO <sub>2</sub>	H	H	H
	1-125	H	NO <sub>2</sub>	H	H	H
35	1-126	4-SO <sub>2</sub> HC <sub>6</sub> H <sub>4</sub>	H	H	H	H
	1-127	Me	SO <sub>2</sub> H	H	H	Et
	1-128	Me	SO <sub>2</sub> H	H	H	H
40	1-129	Me	SO <sub>2</sub> NHMe	H	H	Et
	1-130	Me	SO <sub>2</sub> NHMe	H	H	H
45						
	1-133	Cl	COOEt	H	H	Et
50						
55						

Table 4 (continued)

Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
1-135	Cl	H	H	H	H
1-136	Me	NH <sub>2</sub>	H	H	Et
1-137	Me	C <sub>6</sub> H <sub>5</sub> CH=N	H	H	H
1-138	C <sub>6</sub> H <sub>5</sub> O	H	H	H	H
1-139	2-FC <sub>6</sub> H <sub>4</sub>	H	H	H	Et
1-140	4-FC <sub>6</sub> H <sub>4</sub>	H	H	H	Et
1-141	4-FC <sub>6</sub> H <sub>4</sub>	H	H	H	H
1-142	2,4-F <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	Et
1-143	2,4-F <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	H
1-144	2,6-F <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	Et
1-145	2,6-F <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	H
1-146	Me	NO <sub>2</sub>	H	H	Et
1-147	2-thiazolyl	H	H	H	Et
1-148	2-SMe,5-MeC <sub>6</sub> H <sub>3</sub>	H	H	H	H
1-149	2-Me,4-MeSC <sub>6</sub> H <sub>3</sub>	H	H	H	Et
1-150	4-FC <sub>6</sub> H <sub>4</sub>	H	H	H	H
1-151	2,4-(i-Pr) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	Et
1-152	2,4-(i-Pr) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	H
1-153	2-phenyl-imidazo[1,2-b]pyridazine-8-carboxylic acid				
1-154	6-chloro-2-phenyl-imidazo[1,2-b]pyridazine-8-carboxylic acid				

Table 5

Compound No.	m. p. (°C)	NMR ( $\delta$ value)
1-1	67~68	4.13(3H, s), 6.90(1H, t, J=6.75Hz), 7.70(1H, d, J=0.9Hz), 7.80(1H, d, J=0.9Hz), 8.37(1H, dd, J=6.75Hz, 1.2Hz). (CDCl <sub>3</sub> )
1-2	210 (decomp.)	7.70(1H, d, J=7.5Hz), 8.20(1H, d, J=1.5Hz), 8.55(1H, d, J=7.5Hz), 8.65(1H, d, J=1.5Hz), 9.32(1H, d, J=7.5Hz). (DMSO-d <sub>6</sub> )
1-3	oil	2.45(3H, s), 4.03(3H, s), 6.82(1H, t, J=6.75Hz), 7.45(1H, s), 7.95(1H, d, J=6.75Hz), 8.25(1H, d, J=6.75Hz). (CDCl <sub>3</sub> )
1-4		2.50(3H, s), 7.45(1H, t, J=6.75Hz), 8.15(1H, s), 8.25(1H, d, J=6.75Hz), 9.00(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-5	132~134 (NBr salt)	1.32(3H, t, J=7.0Hz), 2.95(2H, q, J=7.0Hz), 4.05(3H, s), 7.60(1H, d, J=7.5Hz), 8.35(1H, s), 8.50(1H, d, J=6.75Hz), 9.25(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-6	125~127	1.38(3H, t, J=7.5Hz), 3.00(2H, q, J=7.5Hz), 7.62(1H, t, J=6.75Hz), 8.50(1H, d, J=6.75Hz), 8.55(1H, s), 9.45(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-8	oil	1.40(6H, d, J=7.5Hz), 3.00-3.50(1H, m), 4.00(3H, s), 6.75(1H, t, J=6.75Hz), 7.45(1H, s), 7.90(1H, d, J=6.75Hz), 8.26(1H, d, J=6.75Hz, 1.5Hz). (DMSO-d <sub>6</sub> )
1-9	135~137	1.30(6H, d, J=7.5Hz), 3.00-3.60(1H, m), 7.07(1H, t, J=6.75Hz), 7.85(1H, s), 7.99(1H, d, J=6.75Hz), 8.75(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m. p. (°C)	NMR ( $\delta$ value)
1-11	157~158 (HBr salt)	1.48(9H, s), 4.03(3H, s), 7.58(1H, t, J=6.75Hz), 8.33(1H, s), 8.45(1H, d, J=6.75Hz), 9.12(1H, t, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-12	228	1.50(9H, s), 7.69(1H, t, J=6.3Hz), 8.40-8.60(2H, m), 9.26(1H, d, J=6.3Hz). (DMSO-d <sub>6</sub> )
1-13	oil	1.45(3H, t, J=6.0Hz), 4.55(2H, q, J=6.0Hz), 7.05(1H, t, J=6.75Hz), 7.70(1H, s), 8.00(1H, d, J=6.75Hz), 8.28(1H, d, J=6.75Hz). (CDCl <sub>3</sub> )
1-14		7.56(1H, t, J=6.75Hz), 8.20(1H, s), 8.38(1H, d, J=6.75Hz), 8.90(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-15	oil	2.50(3H, s), 4.05(3H, s), 6.95(1H, t, J=6.75Hz), 8.00(1H, d, J=6.75Hz), 9.00(1H, d, J=6.75Hz). (CDCl <sub>3</sub> )
1-17	123~124	2.20(6H, s), 2.50(3H, s), 3.65(2H, s), 6.73(1H, t, J=6.75Hz), 7.95(1H, d, J=6.75Hz), 8.42(1H, d, J=6.75Hz). (CDCl <sub>3</sub> )
1-22	255(sub-limed)	3.20(3H, s), 8.04(1H, t, J=6.75Hz), 9.09(1H, d, J=6.75Hz), 10.04(1H, d, J=6.75Hz). (CF <sub>3</sub> COOD)
1-24	oil	4.05(3H, s), 6.83(1H, t, J=7.5Hz), 7.35-7.60(3H, m), 7.95(1H, s), 7.98-8.15(3H, m), 8.30(1H, d, J=7.5Hz). (CDCl <sub>3</sub> )
1-26	226	7.40-7.70(4H, m), 7.95-8.10(2H, m), 8.33(1H, d, J=6.75Hz), 8.80(1H, s), 9.10(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-27	oil	2.60(3H, s), 4.05(3H, s), 6.88(1H, t, J=6.75Hz), 7.25-7.40(3H, m), 7.80(1H, s), 7.90-8.10(1H, m), 8.35(1H, d, J=6.75Hz). (CDCl <sub>3</sub> )



Table 5 (continued)

Compound No.	m. p. (°C)	NMR ( $\delta$ value)
1-28	220(sub-limed)	2.38(3H, s), 7.30-7.80(5H, m), 8.50(1H, J=6.75Hz), 8.75(1H, s), 9.28(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-30		2.48(3H, s), 4.03(3H, s), 7.30-7.90(5H, m), 8.42(1H, d, J=6.75Hz), 8.87(1H, s), 9.15(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-31	225~227	2.40(3H, s), 7.10-7.50(3H, m), 7.70-7.90(2H, m), 8.10(1H, d, J=6.75Hz), 8.65(1H, s), 8.90(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-32	122~124	2.35(3H, s), 4.06(3H, s), 6.80(1H, d, J=6.75Hz), 7.25(2H, d, J=8.25Hz), 7.85-8.05(3H, m), 7.90(1H, s), 8.28(1H, dd, J=6.75Hz, 1.5Hz). (DMSO-d <sub>6</sub> )
1-33	178~180	2.38(3H, s), 7.30-7.50(3H, m), 7.85(2H, d, J=8.25Hz), 8.25(1H, d, J=6.75Hz), 8.70(1H, s), 9.04(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-34	233~235	2.33(3H, s), 2.48(3H, s), 7.02-7.30(2H, m), 7.67(1H, m), 8.02(1H, dd, J=7.5Hz, 1.5Hz), 8.37(1H, s), 8.85(1H, dd, J=7.5Hz, 1.5Hz). (DMSO-d <sub>6</sub> )
1-36	(HBr salt)	4.00(3H, s), 4.08(3H, s), 7.10-7.40(2H, m), 7.50-7.80(2H, m), 7.93(1H, dd, J=1.5Hz, 7.5Hz), 8.54(1H, d, J=6.75Hz), 8.95(1H, s), 9.34(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-37	238~240	4.00(3H, s), 7.00-7.60(4H, m), 8.05(1H, dd, J=1.5Hz, 7.5Hz), 8.27(1H, d, J=6.75Hz), 8.75(1H, s), 9.10(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m. p. (°C)	NMR ( $\delta$ value)
1-38	172~174 (HBr salt)	4.00(3H, s), 4.17(3H, s), 7.20(1H, m), 7.50-7.80(4H, m), 8.54(1H, d, J=6.75Hz), 8.99(1H, s), 9.25(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-39	209~211	3.90(3H, s), 7.10(1H, m), 7.30-7.75(4H, m), 8.48(1H, d, J=6.75Hz), 9.00(1H, s), 9.25(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-40	oil	3.84(3H, s), 4.05(3H, s), 6.83(1H, d, J= 6.75Hz), 7.00(3H, d, J=8.25Hz), 7.80- 8.10(4H, m), 8.25(1H, dd, J=6.75Hz, 1.5Hz). (CDCl <sub>3</sub> )
1-41	171~173	3.78(3H, s), 7.05(1H, d, J=8.25Hz), 7.30(3H, t, J=6.75Hz), 7.88(2H, d, J= 6.75Hz), 8.60(1H, s), 8.98(1H, d, J= 6.75Hz). (DMSO-d <sub>6</sub> )
1-42	224~225 (HBr salt)	1.43(3H, t, J=6.6Hz), 3.88(3H, s), 3.90(3H, s), 4.47(2H, q, J=6.6Hz), 7.15(1H, d, J=9.0Hz), 7.48-7.68(2H, m), 8.42(1H, dd, J=7.5Hz, 1.5Hz), 8.88(1H, s) 9.12(1H, d, J=7.5Hz). (DMSO-d <sub>6</sub> )
1-43	243~245	3.85(3H, s), 3.90(3H, s), 7.10(1H, d, J= 9.0Hz), 7.40-7.78(2H, m), 8.35(1H, d, J= 7.5Hz), 8.83(1H, s), 9.10(1H, d, J=7.5Hz) (DMSO-d <sub>6</sub> )
1-45	205(sub- limed)	7.25(1H, t, J=6.3Hz), 7.40-7.60(3H, m), 8.00-8.30(2H, m), 8.84(1H, s), 9.00(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )
1-46	220(sub- limed) (HBr salt)	4.05(3H, s), 7.50-7.80(3H, m), 7.95(1H, s), 8.10(1H, br), 8.45(1H, d, J=6.75Hz) 8.96(1H, s), 9.15(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m. p. (°C)	NMR ( $\delta$ value)
1-47	225(sub-limed)	7.40-7.60(3H, m), 7.95(1H, m), 8.15(1H, br), 8.35(1H, d, J=6.75Hz), 8.92(1H, s), 9.11(1H, d, J=6.75Hz), (DMSO-d <sub>6</sub> )
1-48	132~133	4.05(3H, s), 6.85(1H, t, J=6.7Hz), 7.45(2H, d, J=8.3Hz), 7.90(1H, s), 7.80-8.10(3H, m), 8.30(1H, dd, J=6.75Hz, 1.5Hz), (CDCl <sub>3</sub> )
1-49	220~222	7.50-7.70(3H, m), 8.05(2H, d, J=8.3Hz), 8.50(1H, d, J=6.75Hz), 8.85(1H, s), 9.15(1H, d, J=6.75Hz), (DMSO-d <sub>6</sub> )
1-50	160~162 (HBr salt)	4.20(3H, s), 7.60-8.00(5H, m), 8.60-8.80(2H, m), 9.10(1H, d, J=7.78Hz), (DMSO-d <sub>6</sub> )
1-51	205(sub-limed)	7.41(1H, t, J=6.75Hz), 7.68(2H, d, J=8.25Hz), 7.94(2H, d, J=8.25Hz), 8.23(1H, d, J=6.75Hz), 8.85(1H, s), 9.08(1H, d, J=6.75Hz), (DMSO-d <sub>6</sub> )
1-52	157~158	4.00(3H, s), 7.08(1H, t, J=6.75Hz), 7.96(1H, d, J=6.75Hz), 8.35(4H, s), 8.78(1H, s), 8.85(1H, d, J=6.75Hz), (DMSO-d <sub>6</sub> )
1-53	240(sub-limed)	7.28(1H, t, J=6.75Hz), 8.20(1H, d, J=6.75Hz), 8.30-8.50(4H, m), 8.82(1H, s), 8.95(1H, d, J=6.75Hz), (DMSO-d <sub>6</sub> )
1-54	266	7.20(1H, t, J=6.75Hz), 7.25-7.55(3H, m), 8.10(1H, d, J=6.75Hz), 8.18-8.40(1H, m), 8.54(1H, d, J=4.5Hz), 8.95(1H, d, J=6.75Hz), (DMSO-d <sub>6</sub> )
1-56	199~201	2.48(6H, m), 7.15-7.70(4H, m), 8.32(1H, d, J=6.75Hz), 8.55(1H, s), 9.04(1H, d, J=6.75Hz), (DMSO-d <sub>6</sub> )
1-60	97~98	4.18(3H, s), 6.99(1H, t, J=6.9Hz), 7.25-7.63(3H, m), 7.93-8.38(4H, m), (CDCl <sub>3</sub> )
1-61	194~196	7.26(1H, t, J=7.25Hz), 7.92-8.25(3H, m), 8.64(1H, dd, J=1.5Hz, 7.2Hz), (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
1-63	129~131	4.08(3H, s), 6.80(3H, t, J=6.3Hz), 7.21-7.35(3H, m), 7.35-7.62(4H, m), 7.65-7.82(3H, m), 7.98(1H, d, J=6.3Hz), 8.08(1H, d, J=6.3Hz). (CDCl <sub>3</sub> )
1-71	oil	1.43(3H, t, J=6.0Hz), 4.48(2H, q, J=6.0Hz), 7.00(1H, t, J=6.3Hz), 8.01(1H, dd, J=7.5Hz, 0.8Hz), 8.42(1H, dd, J=7.5Hz, 0.8Hz), 8.00-8.10(1H, m). (CDCl <sub>3</sub> )
1-72	203~205	7.20(1H, t, J=6.0Hz), 8.55(1H, dd, J=9.0Hz, 1.5Hz), 8.60-8.67(1H, m), 8.83(1H, dd, J=9.0Hz, 1.5Hz). (DMSO-d <sub>6</sub> )
1-79	89~91	6.71-6.70(2H, m), 6.70-6.95(2H, m), 7.21-7.35(3H, m), 7.35-7.62(4H, m), 7.65-7.82(3H, m), 8.08(1H, d, J=6.3Hz). (DMSO-d <sub>6</sub> )
1-81	250(sub-limed) (NBr salt)	2.50(3H, s), 4.05(3H, s), 7.48-7.70(3H, m), 7.88-8.08(2H, m), 8.40(1H, d, J=1.0Hz), 8.80(1H, s), 9.02(1H, d, J=1.0Hz). (DMSO-d <sub>6</sub> )
1-82	250~252	2.40(3H, s), 7.39-7.65(3H, m), 7.95-8.10(2H, m), 8.09(1H, d, J=1.5Hz), 8.50(1H, s), 8.78(1H, d, J=1.5Hz). (DMSO-d <sub>6</sub> )
1-83	157~158	1.50(3H, t, J=6.3Hz), 2.41(3H, s), 2.55(3H, s), 4.58(2H, q, J=6.3Hz), 6.42(1H, s), 7.20-7.50(4H, m), 7.65(1H, s), 7.85-8.10(2H, m). (CDCl <sub>3</sub> )

Table 5 (continued)

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
1-84	121~122	2.68(3H,s), 2.71(3H,s), 7.05(1H,s), 7.40-7.70(4H,m), 8.60(1H,s). (DMSO-d <sub>6</sub> )
1-85		2.18(3H,s), 3.94(2H,s), 4.02(3H,s), 6.81(1H,t, J=6.6Hz), 7.68(1H,s), 7.93(1H,dd, J=6.6Hz, 1.5Hz), 8.28(1H, dd, J=6.6Hz, 1.5Hz). (DMSO-d <sub>6</sub> )
1-86	250(sub- limed)	5.38(2H,s), 6.88-7.20(3H,m), 7.21-7.5 5(3H,m), 8.27(1H,d, J=7.5Hz), 8.42(1H, s), 9.05(1H,d, J=7.5Hz). (DMSO-d <sub>6</sub> )
1-88	255(sub- limed)	4.48(2H,s), 7.15-7.55(6H,m), 8.13(1H, s), 8.10(1H,d, J=6.6Hz), 8.89(1H,d, J= 6.6Hz). (DMSO-d <sub>6</sub> )
1-89	138~139	1.42(3H,t, J=6.3Hz), 4.47(2H,q, J= 6.3Hz), 4.86(2H,s), 6.87(1H,t, J=7.5Hz) , 7.72(1H,s), 7.93(1H,dd, J=7.5Hz, 1.5Hz), 8.28(1H,dd, J=7.5Hz, 1.5Hz). (CDCl <sub>3</sub> )
1-90	250 (decomp.)	4.86(2H,s), 6.87(1H,t, J=7.5Hz), 7.72(1H,s), 7.93(1H,dd, J=7.5Hz, 1.5Hz) , 8.28(1H,dd, J=7.5Hz, 1.5Hz). (CDCl <sub>3</sub> )
1-94	oil	2.18(3H,s), 3.94(2H,s), 4.02(3H,s), 6.81(1H,t, J=6.6Hz), 7.68(1H,s), 7.93(1H,dd, J=6.6Hz, 1.5Hz), 8.28(1H, dd, J=6.6Hz, 1.5Hz). (CDCl <sub>3</sub> )

Table 5 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
I-96	more than 300	7.60(III,t,J=7.8Hz), 8.04-8.52(5H,m), 9.56-9.72(III,m). (DMSO-d <sub>6</sub> )
I-98	210(sub-limed)	7.22(III,t,J=6.75Hz), 7.83(III,d,J=3.0Hz), 8.00(III,d,J=3.0Hz), 8.05(III,d,J=6.75Hz), 8.67(III,s), 8.92(III,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
I-100	oil	1.32(9H,s), 1.45(3H,t,J=7.0Hz), 2.60(3H,s), 4.52(2H,q,J=7.0Hz), 6.75(III,t,J=6.75Hz), 7.20-7.40(2H,m), 7.70-8.00(3H,m), 8.33(III,d,J=6.75Hz). (CDCl <sub>3</sub> )
I-101	191~192	1.30(9H,s), 2.56(3H,s), 7.20(III,t,J=6.75Hz), 7.36(2H,m), 7.82(III,d,J=9.0Hz), 8.40(III,s), 8.94(III,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
I-102	oil	1.35(9H,s), 1.50(3H,t,J=7.0Hz), 2.51(3H,s), 4.51(2H,q,J=7.0Hz), 6.78(III,t,J=6.75Hz), 7.05-7.40(2H,m), 7.76(III,s), 7.90(III,d,J=6.75Hz), 8.05(III,m), 8.32(III,d,J=6.75Hz). (CDCl <sub>3</sub> )
I-103	194~195	1.32(9H,s), 2.45(3H,s), 7.15(III,t,J=6.75Hz), 7.30(2H,m), 7.85(III,m), 8.08(III,d,J=6.75Hz), 8.45(III,s), 8.90(III,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
I-104	240~241	2.37(3H,s), 2.45(3H,s), 7.17(III,d,J=7.5Hz), 7.23(III,s), 7.60(III,d,J=6.6Hz), 7.45(III,t,J=6.6Hz), 8.28(III,dd,J=6.6,1.5Hz), 8.52(III,s), 9.05(III,dd,J=6.6,1.5Hz). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
I-105	249~250	2.30(3H,s), 2.34(3H,s), 7.31(1H,d, J=6.6Hz), 7.42(1H,d, J=7.5Hz), 7.72(1H,dd, J=7.5Hz, 1.5Hz), 7.77(1H,d, J=1.5Hz), 8.28(1H,dd, J=6.6, 1.5Hz), 8.70(1H,s), 9.03(1H,dd, J=6.6, 1.5Hz). (DMSO-d <sub>6</sub> )
I-106	211~212 (HBr salt)	1.45(3H,t, J=7.0Hz), 2.20-2.45(2H,br), 4.55(2H,q, J=7.0Hz), 7.19(1H,s), 7.63(1H,d, J=6.75Hz), 8.52(1H,d, J=6.75Hz), 8.52(1H,d, J=6.75Hz), 8.59(1H,s), 9.33(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )
I-107	265~266	2.20-2.45(2H,m), 7.15(1H,s), 7.67(1H,d, J=6.75Hz), 8.45-8.60(2H,m), 9.29(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )
I-110	212~213 (HBr salt)	1.28(3H,t, J=7.0Hz), 1.48(3H,t, J=7.0Hz), 2.74(2H,q, J=7.0Hz), 4.57(2H,q, J=7.0Hz), 7.43(2H,d, J=7.5Hz), 7.64(1H,t, J=6.75Hz), 7.95(2H,d, J=7.5Hz), 8.53(1H,d, J=6.75Hz), 8.88(1H,s), 9.25(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )
I-111	214~215	1.25(3H,t, J=7.0Hz), 2.75(2H,q, J=7.0Hz), 7.25-7.60(3H,m), 7.91(2H,d, J=7.5Hz), 8.33(1H,d, J=6.75Hz), 8.79(1H,s), 9.10(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m. p. (°C)	NMR ( $\delta$ value)
1-112	260(sub-limed)	2.35(3H, s), 3.98(3H, s), 7.15(1H, t, J=7.5Hz), 7.30(1H, dd, J=8.0, 1.5Hz), 7.45(1H, d, J=7.5Hz), 7.87(1H, d, J=1.5Hz), 8.25(1H, d, J=7.5Hz), 8.73(1H, s), 9.05(1H, d, J=7.5Hz). (DMSO-d <sub>6</sub> )
1-113	225(sub-limed)	2.37(3H, s), 2.45(3H, s), 7.17(1H, d, J=7.5Hz), 7.23(1H, s), 7.60(1H, d, J=6.6Hz), 7.45(1H, t, J=6.6Hz), 8.28(1H, dd, J=6.6, 1.5Hz), 8.52(1H, s), 9.05(1H, dd, J=6.6, 1.5Hz). (DMSO-d <sub>6</sub> )
1-114	283~285	4.00(3H, s), 7.33(1H, t, J=6.6Hz), 7.49(1H, d, J=6.3Hz), 7.69(1H, dd, J=6.3, 2.5Hz), 8.12(1H, d, J=2.5Hz), 8.40(1H, d, J=6.3Hz), 8.89(1H, s), 9.16(1H, d, J=6.3Hz). (DMSO-d <sub>6</sub> )
1-115	235~237	2.35(3H, s), 3.98(3H, s), 7.15(1H, t, J=7.5Hz), 7.30(1H, dd, J=8.0, 1.5Hz), 7.45(1H, d, J=7.5Hz), 7.87(1H, d, J=1.5Hz), 8.25(1H, d, J=7.5Hz), 8.73(1H, s), 9.05(1H, d, J=7.5Hz). (DMSO-d <sub>6</sub> )
1-118	164~167 (HBr salt)	1.45(3H, t, J=7.0Hz), 2.44(3H, s), 2.53(3H, s), 4.55(2H, q, J=7.0Hz), 7.15-7.40(2H, m), 7.50-7.75(2H, m), 8.47(1H, d, J=6.75Hz), 8.64(1H, s), 9.22(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )



Table 5 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
1-119	139	1.48(3H,t,J=6.9Hz), 4.53(2H,q,J=6.9Hz), 7.20(1H,t,J=7.2Hz), 7.50-7.65(3H,m), 7.82-8.03(2H,m), 8.27(1H,dd,J=1.35, 7.2 Hz), 9.90(1H,dd,J=1.35, 7.2Hz), 10.15 (1H,s). (CDCl <sub>3</sub> )
1-120	214~216	7.35-7.70(4H,m), 7.88-8.09(2H,m), 8.32(1 H,dd,J=1.35, 6.6Hz), 9.80(1H,dd,J=1.35, 6.6Hz), 10.12(1H,s). (DMSO-d <sub>6</sub> )
1-121	87~89	1.49(3H,t,J=7.5Hz), 4.52(2H,q,J=7.5Hz), 6.99(1H,t,J=7.35Hz), 7.36-7.69(3H,m), 8.01(1H,dd,J=1.35, 7.35Hz), 8.17-8.49(3 H,m). (CDCl <sub>3</sub> )
1-122	200~201	7.28(1H,t,J=6.9Hz), 7.40-7.70(3H,m), 8.00-8.23(3H,m), 8.68(1H,dd,J=1.35, 6.9 Hz). (DMSO-d <sub>6</sub> )
1-123	97	1.43(3H,t,J=3.9Hz), 1.45(3H,t,J=3.9Hz), 4.03-4.68(4H,m), 6.99(1H,t,J=6.9Hz), 7. 94-8.12(1H,m), 8.30(1H,s), 8.40(1H,d,J= 6.9Hz). (CDCl <sub>3</sub> )
1-124	139~140	1.39(6H,t,J=7.8Hz), 4.45(2H,q,J=7.8Hz), 4.48(2H,q,J=7.8Hz), 7.63(1H,t,J=7.65Hz), 8.37(1H,dd,J=1.35, 7.65Hz), 9.52(1H,dd, J=1.35, 7.65Hz). (DMSO-d <sub>6</sub> )
1-125	> 300	7.67(1H,t,J=7.5Hz), 8.30(1H,dd,J=1.35, 7.5Hz), 8.80(1H,s), 9.55(1H,dd,J=1.35, 7.5Hz). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m. p. (°C)	NMR ( $\delta$ value)
1-127	247~250 (decomp.)	1.40(3H, t, J=7.5Hz), 2.71(3H, s), 2.90- 4.24(1H, br), 4.51(2H, q, J=7.5 Hz), 7.72 (1H, t, J=7.8Hz), 8.57(1H, dd, J=1.35, 7.8 Hz), 9.30(1H, dd, J=1.35, 7.8Hz). (DMSO-d <sub>6</sub> )
1-128	> 300	2.71(3H, s), 7.70(1H, t, J=7.5Hz), 8.52 (1H, dd, J=1.35, 7.5Hz), 9.30(1H, dd, J= 1.35, 7.5Hz). (DMSO-d <sub>6</sub> )
1-129	184~185	1.42(3H, t, J=7.5Hz), 2.65(3H, s), 2.77(3H, s), 4.47(2H, q, J=7.5Hz), 5.25(1H, br), 7.03 (1H, t, J=7.5Hz), 8.11(1H, dd, J=1.35, 7.5Hz) , 8.87(1H, dd, J=1.35, 7.5Hz). (CDCl <sub>3</sub> )
1-130	235~236	2.49(3H, d, J=5.4Hz), 2.70(3H, s), 5.40- 6.91(1H, br), 7.60(1H, t, J=6.6Hz), 8.34- 8.74(2H, m), 9.02-9.20(1H, m). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m. p. (°C)	NMR ( $\delta$ value)
1-133	105~106	1.44(6H, t, J=7.5Hz), 4.31-4.66(4H, m), 7.15(1H, t, J=7.5Hz), 8.17(1H, dd, J=1.35, 7.5Hz), 9.57(1H, dd, J=1.35, 7.5Hz). (CDCl <sub>3</sub> )
1-135	237	7.25(1H, t, J=7.5Hz), 8.19(1H, dd, J=1.35, 7.5Hz), 7.65-8.29(1H, br), 8.35(1H, s). 8.88(1H, dd, J=1.35, 7.5Hz). (DMSO-d <sub>6</sub> )
1-136	amorphous	1.42(3H, t, J=7.2Hz), 2.43(3H, s), 2.91(2H, br), 4.47(2H, q, J=7.2Hz), 6.81(1H, t, J=7.5 Hz), 7.82(1H, dd, J=1.35, 7.5Hz), 8.20(1H, dd, J=1.35, 7.5Hz). (CDCl <sub>3</sub> )
1-137	231~233	2.70(3H, s), 7.20(3H, t, J=6.9Hz), 7.47- 7.68(3H, m), 7.93-8.21(3H, m), 8.76-8.95 (2H, m). (DMSO-d <sub>6</sub> )
1-139	270(decomp.) (HBr salt)	1.42(3H, t, J=7.0Hz), 4.54(2H, q, J=7.0Hz), 7.30-7.70(4H, m), 8.00-8.25(1H, m), 8.39 (1H, d, J=6.75Hz), 8.77(1H, d, J=4.5Hz), 9.20(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
1-140	177~178 (HBr salt)	1.42(3H,t,J=7.0Hz), 4.55(2H,q,J=7.0Hz), 7.40(2H,t,J=9.0Hz), 7.62(1H,t,J=6.75Hz), 8.03(2H,dd,J=9.0, 6.0Hz), 8.48(1H,d,J=6.75Hz), 8.88(1H,s), 9.20(1H,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
1-141	220 (sublimed)	7.20-7.50(3H,m), 8.05(2H,dd,J=9.0, 6.0Hz), 8.28(1H,d,J=6.75Hz), 8.73(1H,s), 9.06(1H,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
1-142	97~99	1.50(3H,t,J=7.0Hz), 4.65(2H,q,J=7.0Hz), 6.80-7.20(3H,m), 7.90(1H,d,J=6.75Hz), 8.06(1H,d,J=4.0Hz), 8.34(1H,d,J=6.75Hz), 8.45-8.65(1H,m)(CDCl <sub>3</sub> )
1-143	240 (sublimed)	7.00-7.40(2H,m), 7.65-8.10(2H,m), 8.56 (1H,s), 8.85(1H,d,J=6.75Hz), 9.01(1H,d, J=6.75Hz). (CF <sub>3</sub> COOD)
1-144	206~208 (HBr salt)	1.41(3H,t,J=7.0Hz), 4.55(2H,q,J=7.0Hz), 7.20-7.90(4H,m), 8.55(1H,d,J=6.75Hz), 8.88(1H,s), 8.32(1H,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
1-145	280~282	7.15-7.80(3H,m), 8.12(1H,d,J=6.75Hz), 8.55(1H,d,J=1.5Hz), 8.99(1H,d,J=6.75Hz). (DMSO-d <sub>6</sub> )

Table 5 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
1-146	143~144	1.45(3H,t,J=7.0Hz), 2.78(3H,s), 4.42(2H,q,J=7.0Hz), 7.50(1H,d,J=6.75Hz), 8.25(1H,d,J=6.75Hz), 9.55(1H,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
1-147	143~146 (HBr salt)	1.38(3H,t,J=7.0Hz), 4.42(2H,q,J=7.0Hz), 7.30(1H,m), 8.88(2H,m), 8.30(1H,m), 8.70(2H,m). (DMSO-d <sub>6</sub> )
1-151	oil	1.30(2H,d,J=6.0Hz), 1.45(3H,t,J=7.0Hz), 2.95(1H,m,J=6.0Hz), 3.48(1H,m,J=6.0Hz), 4.45(2H,q,J=7.0Hz), 6.75-7.75(5H,m), 7.95(1H,d,J=6.75Hz), 8.35(1H,d,J=6.75Hz). (CDCl <sub>3</sub> )
1-152	99~101	1.15-1.40(12H,m), 2.80-3.40(2H,m), 7.10-7.60(4H,m), 8.35(1H,d,J=6.75Hz), 8.48(1H,s), 9.09(1H,d,J=6.75Hz). (DMSO-d <sub>6</sub> )

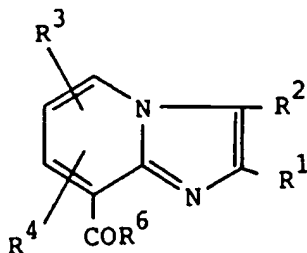



Table 6

Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>a</sup>
5	2-1	H	H	H	A-1
	2-2	Me	H	H	A-1
	2-3	Me	H	H	A-2
10	2-4	Me	H	H	A-4
	2-5	Et	H	H	A-1
	2-6	nPr	H	H	A-1
15	2-7	iPr	H	H	A-1
	2-8	nBu	H	H	A-1
	2-9	tBu	H	H	A-1
20	2-10	H	Cl	H	A-1
	2-11	Me	Cl	H	A-1
	2-12	Me	Me	H	A-1
25	2-13	Me	CHO	H	A-1
	2-14	Me	NO <sub>2</sub>	H	A-1
	2-15	Me	CH <sub>2</sub> NMe <sub>2</sub>	H	A-1
	2-16	Me	CF <sub>2</sub> CHF <sub>2</sub>	H	A-1
30	2-17	Me	CH <sub>2</sub> NMe <sub>2</sub>	H	A-2
	2-18	Me	CH <sub>2</sub> NMe <sub>2</sub>	H	A-4
	2-19	Me	CH <sub>2</sub> NMe <sub>2</sub> · HCl*	H	A-4
35	2-20	Me		H	A-1
	2-21	Me	CH <sub>2</sub> S-iPr	H	A-4
	2-22	Me	CH <sub>2</sub> CH <sub>2</sub> CN	H	A-4
40	2-23	Me	CH=CMe <sub>2</sub>	H	A-4
	2-24	C <sub>6</sub> H <sub>5</sub>	H	H	A-1
	2-25	C <sub>6</sub> H <sub>5</sub>	H	H	A-2
45	2-26	C <sub>6</sub> H <sub>5</sub>	H	H	A-3

T a b l e 6

	Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	2-27	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-4
	2-28	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-5
	2-29	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-6
10	2-30	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-7
	2-31	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-8
	2-32	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-9
15	2-33	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-10
	2-34	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-11
	2-35	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-12
20	2-36	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-13
	2-37	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-14
	2-38	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-15
25	2-39	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-16
	2-40	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-17
	2-41	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-18
30	2-42	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-19
	2-43	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-20
	2-44	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-21
35	2-45	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-22
40	2-48	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-25
	2-49	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-26
45	2-51	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-28
	2-52	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-29

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T a b l e 6

Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>6</sup>
5	2-53	C <sub>6</sub> H <sub>5</sub>	H	H	A-30
	2-54	2-MeC <sub>6</sub> H <sub>4</sub>	H	H	A-1
	2-55	2-MeC <sub>6</sub> H <sub>4</sub>	H	H	A-2
10	2-56	2-MeC <sub>6</sub> H <sub>4</sub>	H	H	A-4
	2-57	3-MeC <sub>6</sub> H <sub>4</sub>	H	H	A-1
	2-58	4-MeC <sub>6</sub> H <sub>4</sub>	H	H	A-1
15	2-59	2,5-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	A-1
	2-60	2,5-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	A-2
	2-61	2,5-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	A-4
20	2-62	2,4,6-Me <sub>3</sub> C <sub>6</sub> H <sub>2</sub>	H	H	A-1
	2-63	2-MeOC <sub>6</sub> H <sub>4</sub>	H	H	A-1
	2-64	3-MeOC <sub>6</sub> H <sub>4</sub>	H	H	A-1
25	2-65	4-MeOC <sub>6</sub> H <sub>4</sub>	H	H	A-1
	2-66	3,4-(MeO) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	A-1
	2-67	3,4-(MeO) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	A-4
30	2-68	2-ClC <sub>6</sub> H <sub>4</sub>	H	H	A-1
	2-69	3-ClC <sub>6</sub> H <sub>4</sub>	H	H	A-1
	2-70	4-ClC <sub>6</sub> H <sub>4</sub>	H	H	A-1
35	2-71	4-BrC <sub>6</sub> H <sub>4</sub>	H	H	A-1
	2-72	4-NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub>	H	H	A-1
	2-73	2-FC <sub>6</sub> H <sub>4</sub>	H	H	A-1
40	2-74	2-F,4-MeC <sub>6</sub> H <sub>3</sub>	H	H	A-1
	2-75	2-Me,4-MeSC <sub>6</sub> H <sub>3</sub>	H	H	A-4
	2-76	2,4,6-Cl <sub>3</sub> C <sub>6</sub> H <sub>2</sub>	H	H	A-1
45	2-77	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> NMe <sub>2</sub>	H	A-1
	2-78	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> NMe <sub>2</sub>	H	A-2

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
Table 6

	Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>a</sup>
5	2-79	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> NMe <sub>2</sub>	H	H	A-4
	2-80	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> NMe <sub>2</sub> · H <sub>2</sub> O*	H	H	A-4
	2-81	C <sub>6</sub> H <sub>5</sub>	Cl	H	H	A-1
10	2-82	C <sub>6</sub> H <sub>5</sub>	Me	H	H	A-4
	2-83	C <sub>6</sub> H <sub>5</sub>	C <sub>6</sub> H <sub>5</sub>	H	H	A-4
	2-84	C <sub>6</sub> H <sub>5</sub>	CF <sub>2</sub> CHF <sub>2</sub>	H	H	A-1
15	2-85	C <sub>6</sub> H <sub>5</sub>	CHO	H	H	A-1
	2-86	C <sub>6</sub> H <sub>5</sub>	NO <sub>2</sub>	H	H	A-1
	2-87	C <sub>6</sub> H <sub>5</sub>	CN	H	H	A-1
20	2-88	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> CH <sub>2</sub> CN	H	H	A-1
	2-89	C <sub>6</sub> H <sub>5</sub>	CH=CMe <sub>2</sub>	H	H	A-1
	2-90	C <sub>6</sub> H <sub>5</sub>	CH(OH)CCl <sub>3</sub>	H	H	A-1
25	2-91	C <sub>6</sub> H <sub>5</sub> CH <sub>2</sub>	H	H	H	A-1
	2-92	CF <sub>3</sub>	H	H	H	A-1
	2-93	Cl	H	H	H	A-1
30	2-94	Cl	Cl	H	H	A-1
	2-95	CHO	H	H	H	A-2
	2-96	CN	H	H	H	A-2
35	2-97	CH=CMe <sub>2</sub>	H	H	H	A-2
	2-98	CH=NOEt	H	H	H	A-2
	2-99	OMe	H	H	H	A-1
40	2-100	OMe	H	H	H	A-2
	2-101	OMe	H	H	H	A-4
	2-102	SMe	H	H	H	A-1
45	2-103	H	OMe	H	H	A-1
	2-104	2-thienyl	H	H	H	A-1
	2-105	2-furyl	H	H	H	A-1

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Table 6

Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>	
5	2-106	2-pyridyl	II	II	II	A-1
	2-107	C <sub>6</sub> H <sub>5</sub>	II	5-Me	II	A-1
	2-108	C <sub>6</sub> H <sub>5</sub>	II	5-Me	II	A-2
10	2-109	C <sub>6</sub> H <sub>5</sub>	II	5-Me	II	A-4
	2-110	Me	II	5-Me	II	A-4
	2-111	Me	CH <sub>2</sub> NMe <sub>2</sub>	5-Me	II	A-2
15	2-112	C <sub>6</sub> H <sub>5</sub>	II	5-Me	7-Me	A-1
	2-113	C <sub>6</sub> H <sub>5</sub>	II	5-CF <sub>3</sub>	7-CF <sub>3</sub>	A-1
	2-114	C <sub>6</sub> H <sub>5</sub> OCH <sub>2</sub>	II	II	II	A-4
20	2-115	C <sub>6</sub> H <sub>5</sub> SCH <sub>2</sub>	II	II	II	A-4
	2-116	C <sub>6</sub> H <sub>5</sub> SOCH <sub>2</sub>	II	II	II	A-4
	2-117	C <sub>6</sub> H <sub>5</sub> SO <sub>2</sub> CH <sub>2</sub>	II	II	II	A-4
25	2-118	C <sub>6</sub> H <sub>5</sub> OCH <sub>2</sub>	CH <sub>2</sub> NMe <sub>2</sub>	II	II	A-4
	2-119	MeOCH <sub>2</sub>	II	II	II	A-1
	2-120	MeOCH <sub>2</sub>	CH <sub>2</sub> N 	II	II	A-1
30	2-121	MeSCH <sub>2</sub>	II	II	II	A-1
	2-122	 MeSCH <sub>2</sub>	II	II	II	A-1
	2-123	MeSCH <sub>2</sub>	CH <sub>2</sub> NMe <sub>2</sub>	II	II	A-1
35	2-124	MeSO <sub>2</sub> CH <sub>2</sub>	II	II	II	A-4
	2-125	 NCH <sub>2</sub>	II	II	II	A-4
	2-126	(EtO) <sub>2</sub> P(O)CH <sub>2</sub>	II	II	II	A-4
40	2-127	Me <sub>2</sub> NCSSCH <sub>2</sub>	II	II	II	A-1
	2-128	ClCH <sub>2</sub>	II	II	II	A-1
	2-129	CF <sub>3</sub>	C <sub>6</sub> H <sub>5</sub> NHCO	II	II	A-4
45	2-130	II	C <sub>6</sub> H <sub>5</sub> CO	II	II	A-2
	2-131	II	C <sub>6</sub> H <sub>5</sub> CH(OH)	II	II	A-2

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Table 6

	Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	2-132	H	C <sub>6</sub> H <sub>5</sub> CH <sub>2</sub>	H	H	A-2
	2-133	H	Et	H	H	A-2
	2-134	C <sub>6</sub> H <sub>5</sub>	H	H	H	A-4
10	2-135	2-MeC <sub>6</sub> H <sub>4</sub>	CHO	H	H	A-1
	2-136	2-FC <sub>6</sub> H <sub>4</sub>	H	H	H	A-4
	2-137	2-FC <sub>6</sub> H <sub>4</sub>	H	H	H	A-2
15	2-138	3,4-MeO <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-139	2,5-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	6-Me	H	A-1
	2-140	4-FC <sub>6</sub> H <sub>4</sub>	H	H	H	A-2
20	2-141	2,4-F <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-142	2,6-F <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-143	C <sub>6</sub> H <sub>5</sub>	CH <sub>2</sub> SO <sub>3</sub> H	H	H	"
25	2-144	4-Me <sub>2</sub> NC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-145	4-NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub>	NO <sub>2</sub>	H	H	"
	2-146	C <sub>6</sub> H <sub>5</sub>	CHO	H	H	"
	2-147	C <sub>6</sub> H <sub>5</sub>	F	H	H	"
30	2-148	C <sub>6</sub> H <sub>5</sub>	CH=NC <sub>6</sub> H <sub>5</sub>	H	H	"
	2-149	4-Me, 3-NO <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	NO <sub>2</sub>	H	H	"
35	2-150	Me	H	H	H	A-11
	2-151	Et	H	H	H	"
	2-152	2-MeC <sub>6</sub> H <sub>4</sub>	H	H	H	"
40	2-153	4-FC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-154	Me	NO <sub>2</sub>	H	H	A-2
	2-155	4-C <sub>6</sub> H <sub>5</sub> -C <sub>6</sub> H <sub>5</sub>	H	H	H	"
45	2-156	C <sub>6</sub> H <sub>5</sub>	H	H	H	"
	2-157	C <sub>6</sub> H <sub>5</sub>	CN	H	H	"

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Table 6

	Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	2-158	H	NO <sub>2</sub>	H	H	"
	2-159	4-Me, 3-NO <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-160	2,4-(NO <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	NO <sub>2</sub>	H	H	"
10	2-161	4-SO <sub>2</sub> HC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-162	2,5-Cl <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-163	2,5-(MeO) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
15	2-164	Me	SO <sub>2</sub> H	H	H	"
	2-165	Me	SO <sub>2</sub> NHMe	H	H	"
	2-166	C <sub>6</sub> H <sub>5</sub>	H	H	H	"
20	2-167	Cl	H	H	H	"
	2-168	4-NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-169	Me	NO <sub>2</sub>	H	H	HNC(OEt)CONH <sub>2</sub>
25	2-170	Me	NO <sub>2</sub>	H	H	A-11
	2-171	2-thiazolyl	H	H	H	A-2
	2-172	CF <sub>3</sub>	H	H	H	HNC(3-FC <sub>6</sub> H <sub>4</sub> )CONH <sub>2</sub>
30	2-173	3-HOC <sub>6</sub> H <sub>4</sub>	H	H	H	A-2
	2-174	2-HOC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-175	2-MeOC <sub>6</sub> H <sub>4</sub>	H	H	H	"
35	2-176	3-MeOC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-177	4-MeOC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-178	2-ClC <sub>6</sub> H <sub>4</sub>	H	H	H	"
40	2-179	3-ClC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-180	4-ClC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-181	2,4-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
45	2-182	3,4-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-183	2-Me, 4-MeSC <sub>6</sub> H <sub>3</sub>	H	H	H	"

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Table 6

Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>5</sup>
5	2-184 2-Me, 4-tBuC <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-185 2-Me, 5-tBuC <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-186 2-Cl, 5-MeC <sub>6</sub> H <sub>3</sub>	H	H	H	"
10	2-187 3-MeC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-188 4-MeC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-189 Me	N=CHC <sub>6</sub> H <sub>5</sub>	H	H	"
15	2-190 OC <sub>6</sub> H <sub>5</sub>	H	H	H	"
	2-191 SC <sub>6</sub> H <sub>5</sub>	H	H	H	"
	2-192 OMe	H	H	H	"
20	2-193 C <sub>6</sub> H <sub>5</sub>	H	5-OH	6-Cl	"
	2-194 $\beta$ -naphthyl	H	H	H	"
	2-195 Me	NH <sub>2</sub>	H	H	"
25	2-196 2,3,4,5-Me <sub>4</sub> C <sub>6</sub> H	H	H	H	"
	2-197 Et	H	H	H	"
	2-198 2-EtC <sub>6</sub> H <sub>4</sub>	H	H	H	"
30	2-199 3-EtC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-200 4-EtC <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-201 4-BrC <sub>6</sub> H <sub>4</sub>	H	H	H	"
35	2-202 2,5-Et <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-203 2-Me, 5-ClC <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-204 2-Me, 5-EtC <sub>6</sub> H <sub>3</sub>	H	H	H	"
40	2-205 2-Et, 5-MeC <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-206 2,4-(MeO) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
	2-207 2-Me <sub>2</sub> NCOC <sub>6</sub> H <sub>4</sub>	H	H	H	"
45	2-208 2-MeOCH <sub>2</sub> C <sub>6</sub> H <sub>4</sub>	H	H	H	"
	2-209 2-Me, 4-MeSOC <sub>6</sub> H <sub>3</sub>	H	H	H	"

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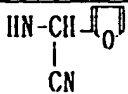
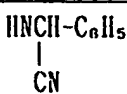


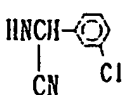
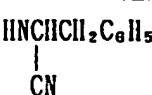
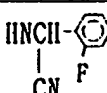

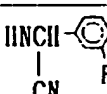
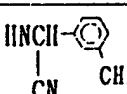
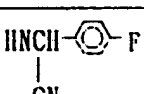
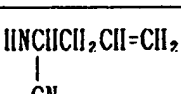
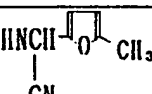
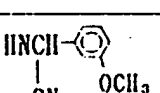
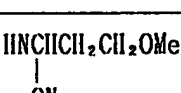

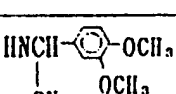


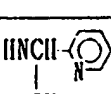
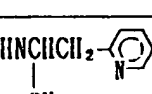
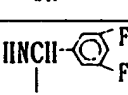
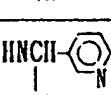

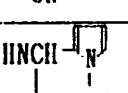
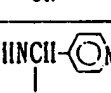
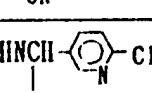
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Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>6</sup>
2-210	2-Me, 4-MeSO <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-211	5-Me, 2-MeOC <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-212	5-Cl, 2-MeOC <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-213	2, 4-(i-Pr) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-214	2, 5-(i-Pr) <sub>2</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-215	4-iPrC <sub>6</sub> H <sub>4</sub>	H	H	H	"
2-216	5-Me, 2-MeSC <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-217	C <sub>6</sub> H <sub>13</sub>	H	H	H	"
2-218	5-t-Bu, 2-MeSC <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-219	5-Cl, 2-MeSC <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-220	4-CNC <sub>6</sub> H <sub>4</sub>	H	H	H	"
2-221	2-CF <sub>3</sub> C <sub>6</sub> H <sub>4</sub>	H	H	H	"
2-222	3-CF <sub>3</sub> C <sub>6</sub> H <sub>4</sub>	H	H	H	"
2-223	2-Me, 5-CF <sub>3</sub> C <sub>6</sub> H <sub>3</sub>	H	H	H	"
2-224	α-(imidazo[1, 2-a]pyridin-5-ylcarbonylamino)-(2-furyl)acetonitrile				
2-225	α-(imidazo[1, 2-a]pyridin-6-ylcarbonylamino)-(2-furyl)acetonitrile				

Compound No.	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>	R <sup>4</sup>	R <sup>6</sup>
2-231	α-(2-phenyl-imidazo[1, 2-b]pyridazin-8-ylcarbonylamino)-(2-thienyl)acetonitrile				
2-232	α-(6-chloro-2-phenyl-imidazo[1, 2-b]pyridazin-8-ylcarbonylamino)-(2-thienyl)acetonitrile				

\*HCl:hydrochloride, \*HlO:oxalate

A-1 to A-30 for R<sup>6</sup> in Table 6 mean the following groups.

5	A-1		A-11		A-21	
10	A-2		A-12		A-22	
15	A-3		A-13			
	A-4		A-14			
20	A-5		A-15		A-25	
25	A-6		A-16		A-26	
30	A-7		A-17			
35	A-8		A-18		A-28	
40	A-9		A-19		A-29	
45	A-10		A-20		A-30	

T a b l e 7

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
2-1	149~151	6.40-6.55(2H, m), 6.65(1H, d, J=3.0Hz), 6.90(1H, t, J=6.75Hz), 7.50(1H, br), 7.55(1H, d, J=0.9Hz), 7.72(1H, d, J=0.9Hz); 8.12(1H, d, J=6.75Hz), 8.32(1H, d, J=6.75Hz), 11.30(1H, d, J=8.75Hz). (CDCl <sub>3</sub> )
2-2	130~131	2.45(3H, s), 6.50(2H, m), 6.67(1H, d, J=3.0Hz), 6.90(1H, t, J=6.75Hz), 7.40(1H, br), 7.55(1H, d, J=1.5Hz), 8.10-8.30(2H, m), 11.45(1H, d, J=7.5Hz). (CDCl <sub>3</sub> )
2-3	140~141	2.40(3H, s), 6.78(1H, d, J=7.5Hz), 7.15(2H, m), 7.45(1H, m), 7.66(1H, d, J=6.0Hz), 7.88(1H, s), 8.08(1H, d, J=6.75Hz), 8.78(1H, d, J=6.75Hz), 11.28(1H, d, J=7.5Hz). (DMSO-d <sub>6</sub> )
2-4	204~205	2.40(3H, s), 6.55(1H, d, J=7.5Hz), 7.08(1H, t, J=6.75Hz), 7.20-7.70(4H, m), 7.88(1H, s), 8.05(1H, d, J=6.75Hz), 8.75(1H, d, J=6.75Hz), 11.24(1H, d, J=7.5Hz). (DMSO-d <sub>6</sub> )
2-5	102~103	1.30(3H, t, J=7.5Hz), 2.80(2H, q, J=7.5Hz), 6.51-6.70(2H, m), 6.75(1H, d, J=3.0Hz), 7.08(1H, t, J=6.75Hz), 7.83(1H, m), 7.88(1H, s), 8.07(1H, dd, J=6.75Hz, 1.5Hz), 8.78(1H, dd, J=6.75Hz, 1.5Hz), 11.35(1H, d, J=6.75Hz). (DMSO-d <sub>6</sub> )



Table 7 (continued)

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
2-7	85~87	1.35(6H, d, J=6.0Hz), 3.08(1H, m), 6.45-6.70(2H, m), 6.70(1H, d, J=3.0Hz), 7.03(1H, t, J=6.75Hz), 7.75(1H, m), 7.80(1H, s), 8.06(1H, dd, J=6.75Hz, 1.5Hz), 8.73(1H, dd, J=6.75Hz, 1.5Hz), 11.48(1H, d, J=6.75Hz). (DMSO <sub>d</sub> )
2-9	112~114	1.32(9H, s), 6.48-6.65(2H, m), 6.73(1H, d, J=3.0Hz), 7.05(1H, t, J=6.75Hz), 7.78(1H, m), 7.84(1H, s), 8.07(1H, d, J=6.75Hz), 8.74(1H, d, J=6.75Hz), 11.53(1H, d, J=6.75Hz). (DMSO <sub>d</sub> )
2-10	142~143	6.50(2H, m), 6.65(1H, d, J=3.75Hz), 7.15(1H, t, J=6.75Hz), 7.45(1H, d, J=1.5Hz), 7.58(1H, s), 8.25(2H, m), 10.95(1H, d, J=7.5Hz). (CDCl <sub>3</sub> )
2-11	157~159	2.45(3H, s), 6.35-6.50(2H, m), 6.65(1H, d, J=3.95Hz), 7.08(1H, t, J=6.75Hz), 7.50(1H, d, J=1.5Hz), 8.10-8.30(2H, m), 11.05(1H, d, J=8.25Hz). (CDCl <sub>3</sub> )
2-15	162~164	2.20(6H, s), 2.40(3H, s), 3.65(2H, s), 6.50(2H, m), 6.68(1H, d, J=3.0Hz), 6.92(1H, t, J=6.75Hz), 7.48(1H, br), 8.12(1H, d, J=6.75Hz), 8.40(1H, d, J=6.75Hz), 11.58(1H, d, J=7.50Hz). (CDCl <sub>3</sub> )
2-17	123~124	2.20(6H, s), 2.40(3H, s), 3.70(2H, s), 6.47(1H, d, J=7.5Hz), 6.95(1H, d, J=6.75Hz), 7.10-7.50(4H, m), 8.20(1H, d, J=6.75Hz), 8.45(1H, d, J=6.75Hz), 11.58(1H, d, J=7.50Hz). (CDCl <sub>3</sub> )
2-19	171~174	

Table 7 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-20	184~186	2.35-2.60(4H,m), 2.42(3H,s), 3.55-3.75(4H,m), 3.77(2H,s), 6.50(2H,m), 6.65(1H,d, J=3.0Hz), 6.95(1H,t, J=6.75Hz), 7.55(1H,br), 8.25(1H,d, J=6.75Hz), 8.47(1H,d, J=6.75Hz), 11.45(1H,d, J =7.5Hz). (DMSO <sub>d</sub> )
2-24	197~199	6.35-6.50(2H,m), 6.68(1H,d, J=3.75Hz), 6.95 (1H,t, J=6.75Hz), 7.35-7.65(4H,m), 7.70-7.90 (2H,m), 7.92(1H,s), 8.10-8.30(2H,m), 11.56(1H, d, J=7.50Hz). (CDCl <sub>3</sub> )
2-25	187~189	6.76(1H,d, J=6.3Hz), 7.05-7.32(2H,m), 7.33- 7.65(4H,m), 7.70-7.82(1H,m), 7.86-8.02(2H,m), 8.04-8.20(1H,m), 8.59(1H,s), 8.73-8.90(1H,m), 11.34(1H,d, J=6.3Hz). (DMSO <sub>d</sub> )
2-26	176~177	6.60(1H,d, J=7.5Hz), 7.14(1H,t, J=6.75Hz), 7.35-7.60(7H,m), 7.80-8.00(2H,m), 8.15(1H,d, J=6.75Hz), 8.67(1H,s), 8.84(1H,d, J=6.75Hz), 11.48(1H,d, J=7.5Hz). (DMSO <sub>d</sub> )
2-27	192~193	6.50(1H,d, J=6.6Hz), 7.14(1H,t, J=7.5Hz), 7.36 -7.73(7H,m), 7.80-7.99(2H,m), 8.02-8.20 (1H,m), 8.56(1H,s), 8.72-8.89(1H,m), 11.30(1H, d, J=6.6Hz). (DMSO <sub>d</sub> )
2-28	205~206	6.36(1H,d, J=6.6Hz), 6.98-7.53(6H,m), 7.68- 7.95(4H,m), 8.03-8.20(1H,m), 8.42(1H,s), 8.66 -8.80(1H,m), 11.33(1H,d, J=6.6Hz). (CDCl <sub>3</sub> + DMSO <sub>d</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR( $\delta$ value)
2-29	168~170	3.82(3H,s), 6.35(1H,d, J=6.3Hz), 7.03-7.66(8H,m), 7.78-7.95(2H,m), 8.10(1H,dd, J=1.35Hz, 7.5Hz), 8.54(1H,s), 8.78(1H,dd, J=1.35Hz, 7.5Hz), 11.28(1H,d, J=6.3Hz). (DMSO-d <sub>6</sub> )
2-30	181~182	3.80(3H,s), 3.85(3H,s), 6.24(1H,d, J=6.3Hz), 7.04-7.53(7H,m), 7.70-7.91(2H,m), 8.10(1H,dd, J=1.5Hz, 7.2Hz), 8.53(1H,s), 8.79(1H,dd, J=1.5Hz, 7.2Hz), 11.22(1H,d, J=6.3Hz). (DMSO-d <sub>6</sub> )
2-31	203~204	6.48(1H,d, J=6.45Hz), 7.12(1H,t, J=7.2Hz), 7.32-7.65(4H,m), 7.68-8.02(5H,m), 8.10(1H,dd, J=1.35, 7.2Hz), 8.55(1H,s), 8.80(1H,dd, J=1.35, 7.2Hz), 11.25(1H,d, J=6.45Hz). (DMSO-d <sub>6</sub> )
2-32	219~220	6.55(1H,d, J=6.75Hz), 7.15(1H,t, J=7.05Hz), 7.24-7.65(6H,m), 7.88-8.03(2H,m), 8.10(1H,dd, J=1.35, 7.05Hz), 8.53(1H,s), 8.82(1H,dd, J=1.35, 7.05Hz), 11.30(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )
2-33	207~208	3.71(3H,s), 6.10-6.26(1H,m), 6.43-6.60(2H,m), 6.89-7.01(1H,m), 7.13(1H,t, J=7.5Hz), 7.31-7.58(3H,m), 7.72-7.92(2H,m), 8.11(1H,dd, J=1.35, 7.5Hz), 8.55(1H,s), 8.79(1H,dd, J=1.35, 7.5Hz), 11.13(1H,d, J=6.45Hz). (DMSO-d <sub>6</sub> )
2-34	174~175	6.28(1H,d, J=7.5Hz), 6.95(1H,t, J=6.9Hz), 7.23-7.84(10H,m), 7.90(1H,s), 8.11-8.32(2H,m), 11.48(1H,d, J=7.5Hz). (CDCl <sub>3</sub> )

Table 7 (continued)

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
2-35	210~211	6.32(1H, d, J=7.5Hz), 7.03(1H, t, J=7.2Hz), 7.17-7.92(4H, m), 8.16(1H, dd, J=1.35Hz, 7.2Hz), 8.22(1H, s), 8.57(1H, dd, J=1.35Hz, 7.2Hz), 11.46(1H, d, J=7.5Hz). (DMSO-d <sub>6</sub> +CDCl <sub>3</sub> )
2-36	192~193	6.52(1H, d, J=7.35Hz), 7.05(1H, t, J=7.2Hz), 7.32-7.53(3H, m), 7.61-7.91(4H, m), 7.92-8.22(3H, m), 8.33(1H, s), 8.65(1H, dd, J=1.35Hz, 7.2Hz), 11.41(1H, d, J=7.35Hz). (CDCl <sub>3</sub> +DMSO-d <sub>6</sub> )
2-37	189~190	2.42(3H, s), 6.34(1H, d, J=6.3Hz), 7.14(1H, t, J=7.2Hz), 7.30-7.68(7H, m), 7.74-7.92(2H, m), 8.11(1H, dd, J=1.35Hz, 7.2Hz), 8.55(1H, s), 8.79(1H, dd, J=1.35Hz, 7.2Hz), 11.25(1H, d, J=6.3Hz). (DMSO-d <sub>6</sub> )
2-40	168~169	1.23(6H, d, J=7.2Hz), 2.12-2.50(1H, m), 5.05(1H, dd, J=6.9Hz, 7.5Hz), 7.16(1H, t, J=6.9Hz), 7.28-7.65(3H, m), 7.90-8.20(3H, m), 8.60(1H, s), 8.82(1H, dd, J=1.35Hz, 6.9Hz), 10.90(1H, d, J=7.5Hz). (DMSO-d <sub>6</sub> )
2-41	181~182	6.37(1H, d, J=7.5Hz), 6.95(1H, t, J=7.2Hz), 6.27-8.37(11H, m), 8.79-8.94(1H, m), 11.97(1H, d, J=7.50Hz). (CDCl <sub>3</sub> )
2-42	213~215	6.58(1H, d, J=6.45Hz), 7.13(1H, t, J=7.2Hz), 7.31-7.71(4H, m), 7.80-8.30(4H, m), 8.55(1H, s), 8.68-8.87(2H, m), 8.96(1H, d, J=2.85Hz), 11.30(1H, d, J=6.45Hz). (CDCl <sub>3</sub> )

Table 7 (Continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-45	182~183	3.31(2H,d,J=7.5Hz),5.32-5.62(1H,m),6.93(1H,t,J=6.6Hz),7.23-7.62(8H,m),7.65-8.02(3H,m),8.07-8.36(2H,m),11.03(1H,d,J=8.7Hz). (CDCl <sub>3</sub> )
2-48	192~193	2.42(3H,s),6.20-6.31(1H,m),6.56(1H,d,J=3.3Hz),7.15(1H,t,J=7.2Hz),7.28-7.63(3H,m),7.89-8.06(3H,m),8.12(1H,d,J=1.35Hz,7.2Hz),8.59(1H,s),8.81(1H,dd,J=1.35Hz,7.2Hz),11.37(1H,d,J=6.9Hz). (DMSO-d <sub>6</sub> )
2-49	173~174	4.60(2H,d,J=6.3Hz),7.12(1H,t,J=6.9Hz),7.24-7.62(3H,m),8.00-8.25(3H,m),8.58(1H,s),8.71-8.87(1H,m),10.59(1H,t,J=6.3Hz). (DMSO-d <sub>6</sub> )

Table 7 (Continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-52	196~197	2.41(3H,s), 6.49(1H,d,J=7.8Hz), 6.80-7.06(2H,m), 7.23-7.57(4H,m), 7.75-7.95(3H,m), 8.07-8.32(2H,m), 11.48(1H,d,J=7.8Hz). (CDCl <sub>3</sub> )
2-53	159~161	6.60(1H,d,J=6.6Hz), 7.12(1H,t,J=6.9Hz), 7.32-7.61(3H,m), 7.69(1H,d,J=9.3Hz), 7.82-8.14(3H,m), 8.22(1H,dd,J=1.2, 9.3Hz), 8.57(1H,s), 8.72-8.89(2H,m), 11.26(1H,d,J=6.6Hz). (DMSO-d <sub>6</sub> )
2-54	175~177	2.54(3H,s), 6.55-6.70(2H,m), 6.75(1H,d,J=3.0Hz), 7.20(1H,t,J=6.75Hz), 7.30-7.50(3H,m), 7.80-8.00(2H,m), 8.15(1H,d,J=6.75Hz), 8.45(1H,s), 8.85(1H,d,J=6.75Hz), 11.35(1H,d,J=7.50Hz). (DMSO-d <sub>6</sub> )
2-55	175~176	2.45(3H,s), 6.75(1H,d,J=7.5Hz), 7.10-8.30(8H,m), 8.42(1H,s), 8.86(1H,d,J=6.75Hz), 11.27(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-56	119~120	2.45(3H,s), 6.53(1H,d,J=7.5Hz), 7.10-7.90(9H,m), 8.15(1H,d,J=6.75Hz), 8.45(1H,s), 8.86(1H,d,J=6.75Hz), 11.28(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-57	214~216	2.40(3H,s), 6.50-6.70(2H,m), 6.82(1H,d,J=3.0Hz), 7.00-7.55(3H,m), 7.65-7.95(3H,m), 8.12(1H,d,J=6.75Hz), 8.55(1H,s), 8.80(1H,d,J=6.75Hz), 11.45(1H,d,J=6.75Hz). (DMSO-d <sub>6</sub> )

Table 7 (Continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-58	198~200	2.35(3H,s), 6.60-6.75(2H,m), 6.80(1H,d, J=3.0Hz), 7.15(1H,t, J=6.75Hz), 7.32(2H,d, J=8.35Hz), 7.75-8.00(2H,m), 8.10(1H,d, J=6.75Hz), 8.54(1H,s), 8.80(1H,d, J=6.75Hz), 11.52(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )
2-59	173~175	2.39(3H,s), 2.50(3H,s), 6.40(1H,d, J=6.6Hz), 6.41-6.58(2H,m), 6.72(1H,d, J=3.0Hz), 7.08(1H,t, J=6.0Hz), 7.18(1H,m), 7.52-7.72(2H,m), 8.05(1H,s), 8.27(1H,dd, J=7.5Hz, 0.3Hz), 8.59(1H,dd, J=7.5Hz, 0.3Hz), 11.48(1H,d, J=6.6Hz). (CDCl <sub>3</sub> )
2-60	134~135	2.34(3H,s), 2.47(3H,s), 6.53(1H,d, J=7.5Hz), 6.90-7.25(4H,m), 7.31-7.48(2H,m), 7.80(1H,s), 7.18-7.40(2H,m), 11.56(1H,d, J=7.5Hz). (CDCl <sub>3</sub> )
2-63	167~169	4.01(3H,s), 6.55-6.70(2H,m), 6.80(1H,d, J=3.0Hz), 7.00-7.50(4H,m), 7.95(1H,br), 8.00-8.25(2H,m), 8.68(1H,s), 8.87(1H,t, J=6.75Hz), 11.51(1H,d, J=6.0Hz). (DMSO-d <sub>6</sub> )
2-64	167~169	3.82(3H,s), 6.55-6.70(2H,m), 6.82(1H,d, J=3.0Hz), 6.85-7.60(5H,m), 7.89(1H,m), 8.13(1H,d, J=6.75Hz), 8.58(1H,s), 8.80(1H,d, J=6.75Hz), 11.42(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )
2-65	179~180	3.85(3H,s), 6.55-6.70(2H,m), 6.70(1H,d, J=3.0Hz), 7.00-7.25(3H,m), 7.85-8.05(3H,m), 8.12(1H,d, J=6.75Hz), 8.45(1H,s), 8.80(1H,d, J=6.75Hz), 11.45(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )

Table 7 (Continued)

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
2-66	186~188	3.84(6H, s), 6.58-6.71(2H, m), 6.80(1H, d, J=3.0 Hz), 7.43-7.68(2H, m), 7.10(1H, d, J=7.5 Hz), 8.10(1H, d, J=6.75 Hz), 8.50(1H, s), 8.80(1H, d, J=6.75 Hz), 11.40(1H, d, J=7.5 Hz). (DMSO-d <sub>6</sub> )
2-67	140~141	3.83(3H, s), 3.90(3H, s), 6.40(1H, d, J=7.5 Hz), 6.83-7.20(3H, m), 7.25-7.60(5H, m), 7.88(1H, s), 8.18(1H, dd, J=6.0 Hz, 1.5 Hz), 8.25(1H, dd, J=6.0 Hz, 1.5 Hz), 11.55(1H, d, J=7.5 Hz). (CDCl <sub>3</sub> )
2-68	176~178	6.55-6.70(2H, m), 6.78(1H, d, J=3.0 Hz), 7.18(1H, t, J=6.75 Hz), 7.34-7.66(3H, m), 7.78(1H, br), 8.10-8.30(2H, m), 8.76(1H, s), 8.88(1H, d, J=6.75 Hz), 11.30(1H, d, J=6.75 Hz). (DMSO-d <sub>6</sub> )
2-69	235~237	6.55-6.75(2H, m), 6.82(1H, d, J=3.0 Hz), 7.20(1H, t, J=6.75 Hz), 7.35-7.70(2H, m), 7.88-8.30(4H, m), 8.66(1H, s), 8.84(1H, d, J=6.75 Hz), 11.32(1H, d, J=6.75 Hz). (DMSO-d <sub>6</sub> )
2-70	222~224	6.55-6.70(2H, m), 6.83(1H, d, J=3.0 Hz), 7.18(1H, t, J=6.75 Hz), 7.53(2H, d, J=8.25 Hz), 7.90-8.20(4H, m), 8.61(1H, s), 8.82(1H, d, J=6.75 Hz), 11.38(1H, d, J=6.75 Hz). (DMSO-d <sub>6</sub> )
2-71	233~234	6.55-6.75(2H, m), 6.82(1H, d, J=3.0 Hz), 7.16(1H, t, J=6.75 Hz), 7.78(2H, d, J=8.25 Hz), 7.90(2H, d, J=8.25 Hz), 7.98(1H, m), 8.15(1H, d, J=6.75 Hz), 8.60(1H, s), 8.83(1H, d, J=6.75 Hz), 11.32(1H, d, J=6.75 Hz). (DMSO-d <sub>6</sub> )



Table 7 (Continued)

Compound No. m.p.(°C)		NMR ( $\delta$ value)
2-72	249~251	6.55-6.70(2H,m), 6.78(1H,d, J=3.0Hz), 7.19 (1H,t, J=6.75Hz), 7.95-8.45(6H,m), 8.75-8.90 (2H,m), 11.25(1H,d, J=6.75Hz). (DMSO-d <sub>6</sub> )
2-73	208~210	6.65(2H,m), 6.83(1H,d, J=3.0Hz), 7.21(1H,t, J=6.75Hz), 7.30-7.60(3H,m), 7.95(1H,m), 8.05-8.35(2H,m), 8.54(1H,d, J=4.5Hz), 8.90(1H,d, J=6.75Hz), 11.46(1H,d, J=7.5Hz). (DMSO-d <sub>6</sub> )
2-78	173~175	2.26(6H,s), 3.92(2H,s), 6.53(1H,d, J=8.1Hz), 6.90-7.11(2H,m), 7.32-7.55(4H,m), 7.63-7.84(2H,m), 8.28(1H,dd, J=7.5Hz, 0.75Hz), 8.63(1H,dd, J=7.5Hz, 0.75Hz), 11.65(1H,d, J=8.1Hz). (CDCl <sub>3</sub> )

Table 7 (continued)

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
2-79	176~177	2.26(6H, s), 3.93(2H, s), 6.36(1H, d, J=8.1Hz), 6.91-7.34(2H, m), 7.43-7.60(6H, m), 7.66-7.85 (2H, m), 8.21-8.38(1H, m), 8.57-8.73(1H, m), 11.65(1H, d, J=8.1Hz). (CDCl <sub>3</sub> )
2-80	~125	
2-81	173	6.58-6.70(2H, m), 6.80(1H, d, J=3.3Hz), 7.32(1H, t, J=6.9Hz), 7.40-7.73(3H, m), 7.88-8.29(4H, m), 8.65(1H, dd, J=1.5Hz, 6.9Hz), 11.05(1H, d, J=7.2 Hz). (DMSO-d <sub>6</sub> )
2-82	169~171	2.70(3H, s), 6.48(1H, d, J=6.3Hz), 7.18(1H, t, J= 6.9Hz), 7.28-7.86(9H, m), 8.12(1H, dd, J=1.5Hz, 6.9Hz), 8.62(1H, dd, J=1.5Hz, 6.9Hz), 11.39(1H, d, J=6.3Hz). (DMSO-d <sub>6</sub> )
2-83	129~131	6.32(1H, d, J=6.3Hz), 6.91(1H, t, J=6.3Hz), 7.15 -7.75(14H, m), 8.05(1H, dd, J=6.3Hz, 1.5Hz), 8.25(1H, dd, J=6.3Hz, 1.5Hz), 11.68(1H, d, J=6.3 Hz). (CDCl <sub>3</sub> )
2-92	145~146	6.31(1H, d, J=8.4Hz), 7.12(1H, t, J=6.5Hz), 7.20 -7.55(4H, m), 7.98(1H, s), 7.30(1H, dd, J=7.5Hz, 1.5Hz), 7.35(1H, dd, J=7.5Hz, 1.5Hz), 10.85(1H, d, J=8.4Hz). (CDCl <sub>3</sub> )
2-104	221~222	6.15-6.70(2H, m), 6.70-6.95(2H, m), 7.20(1H, t, J=6.5Hz), 7.80-7.95(1H, m), 8.27(1H, dd, J=7.5 Hz, 1.5Hz), 8.41(1H, s), 8.83(1H, dd, J=7.5Hz, 1.5Hz), 11.50(1H, d, J=6.0Hz). (DMSO-d <sub>6</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR( $\delta$ value)
2-109	197~198	2.42(3H,s), 6.29(1H,d, J=8.4Hz), 7.15-7.55(7H,m), 7.70-7.85(2H,m), 7.80(1H,s), 8.05(2H,s), 11.53(1H,d, J=8.4Hz). (CDCl <sub>3</sub> )
2-112	206~207	2.57(3H,s), 2.71(3H,s), 6.32(1H,d, J=6.0Hz), 6.50(1H,d, J=6.3Hz), 6.54(1H,d, J=6.3Hz), 6.67(1H,d, J=3.0Hz), 7.31-7.52(3H,m), 7.55(1H,m), 7.62(1H,s), 7.71-7.92(2H,m), 12.30(1H,d, J=6.0Hz). (CDCl <sub>3</sub> )
2-114	119~121	4.22(2H,s), 6.38(1H,d, J=9.0Hz), 6.92(1H,t, J=8.0Hz), 7.10-7.50(9H,m), 7.55(1H,s), 8.19(2H,d, J=6.3Hz), 11.23(1H,d, J=9.0Hz). (CDCl <sub>3</sub> )
2-115	138~139	5.28(2H,s), 6.44(1H,d, J=9.0Hz), 6.95(1H,t, J=8.0Hz), 7.00-7.50(9H,m), 7.70(1H,s), 8.25(2H,d, J=6.3Hz), 11.21(1H,d, J=9.0Hz). (CDCl <sub>3</sub> )
2-119	127~128	3.50(3H,s), 4.65(2H,s), 6.33-6.50(2H,m), 6.93(1H,t, J=6.3Hz), 7.49(1H,m), 7.62(1H,s), 8.15-8.32(2H,m), 11.25(1H,d, J=7.5Hz). (CDCl <sub>3</sub> )
2-120	108~109	2.31-2.57(4H,m), 3.47(3H,s), 3.55-3.75(4H,m), 3.90(2H,s), 4.63(2H,s), 6.37-6.53(2H,m), 6.93(1H,t, J=6.3Hz), 7.48(1H,m), 8.23(1H,dd, J=7.5Hz, 1.5Hz), 8.51(1H,dd, J=7.5Hz, 1.5Hz), 11.32(1H,d, J=7.5Hz). (CDCl <sub>3</sub> )
2-121	165~166	2.17(3H,s), 3.80(2H,s), 6.33-6.50(2H,m), 6.60-6.73(1H,m), 6.95(1H,d, J=6.0Hz), 7.50(1H,m), 7.60(1H,s), 8.12-8.30(2H,m), 11.27(1H,d, J=7.5Hz). (CDCl <sub>3</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR( $\delta$ value)
2-122	214~216	2.61(3H,d,J=3.0Hz), 4.20(2H,s), 6.33-6.51(2H,m), 6.58-6.75(1H,m), 7.08(1H,t, J=6.0Hz), 7.60(1H,m), 8.00(1H,s), 8.18(1H,dd,J=6.0Hz,1.5Hz), 8.60(1H,dd,J=6.0Hz,1.5Hz), 11.08(1H,d,J=6.3Hz). (CDCl <sub>3</sub> +DMSO-d <sub>6</sub> )
2-123	117~118	2.19(3H,s), 2.25(6H,s), 3.72(2H,s), 3.83(2H,s), 6.32-6.48(2H,m), 6.65(1H,d, J=3.0Hz), 6.98(1H,d,J=6.0Hz), 7.50(1H,m), 7.72(1H,dd,J=6.0Hz,1.5Hz), 8.50(1H,dd,J=6.0Hz,1.5Hz), 11.40(1H,d,J=6.6Hz). (CDCl <sub>3</sub> )
2-127	169~171	3.38(3H,br), 3.55(3H,br), 4.75(2H,s), 6.36-6.51(2H,m), 6.59-6.70(1H,m), 6.93(1H,t,J=6.0Hz), 7.45-7.55(1H,m), 7.78(1H,s), 8.28(1H,d,J=7.5Hz), 11.21(1H,br). (CDCl <sub>3</sub> )
2-128	184~186	4.73(2H,s), 6.32-6.50(2H,m), 6.65(1H,d, J=3.0Hz), 6.98(1H,d,J=6.0Hz), 7.50(1H,m), 7.72(1H,s), 8.23(1H,d,J=7.5Hz), 11.10(1H,br). (CDCl <sub>3</sub> )
2-134	120~125 (hydrochloride)	5.70(1H,br), 6.50(1H,d,J=6.3Hz), 7.16-7.82(8H,m), 7.82-8.10(2H,m), 8.16-8.32(1H,m), 8.68(1H,s), 8.82-9.00(1H,m), 11.21(1H,d,J=6.3Hz). (DMSO-d <sub>6</sub> )
2-135	112~115	2.40(3H,s), 6.35(2H,m), 6.63(1H,d,J=3.0Hz), 7.20-7.60(8H,m), 8.55(1H,d,J=6.75Hz), 9.80(1H,d,J=6.75Hz), 9.82(1H,s), 11.07(1H,d,J=8.5Hz). (CDCl <sub>3</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR( $\delta$ value)
2-136	171~172	6.50(1H,d,J=7.5Hz), 7.10-7.75(8H,m), 7.95(1H,m), 8.18(1H,d,J=6.75Hz), 8.50 (1H,d,J=4.5Hz), 8.90(1H,d,J=6.75Hz), 11.20(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-137	190~191	6.75(1H,d,J=7.5Hz), 7.05-7.55(6H,m), 7.72(1H,d,J=6.0Hz), 7.95-8.25(1H,m), 8.15(1H,d,J=6.75Hz), 8.48(1H,d,J= 4.5Hz), 8.88(1H,d,J=6.75Hz), 11.25(1H, d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-138	192~193	3.83(3H,s), 3.91(3H,s), 6.58(1H,d,J= 7.5Hz), 6.80-7.15(3H,m), 7.82(1H,s), 8.23(2H,m), 11.58(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-139	184~185	2.38(6H,s), 2.43(3H,s), 6.32(1H,d,J= 7.5Hz), 6.32-6.49(2H,m), 6.61(1H,m), 6.95-7.20(2H,m), 7.50(1H,m), 6.62(2H, m), 8.05(1H,m), 11.50(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-140	215~217	6.77(1H,d,J=7.5Hz), 7.08-7.45(4H,m), 7.52(1H,d,J=3.0Hz), 7.77(1H,dd,J=6.0, 3.0Hz), 7.85-8.20(3H,m), 8.58(1H,s), 8.80(1H,d,J=6.75Hz), 11.25(1H,d,J= 7.5Hz). (DMSO-d <sub>6</sub> )
2-141	220~222	6.75(1H,d,J=7.5Hz), 7.10-7.50(5H,m), 7.77(1H,dd,J=3.75,5.25Hz), 7.90-8.35 (2H,m), 8.50(1H,d,J=4.0Hz), 8.90(1H,d, J=6.75Hz), 11.13(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-142	154~156	6.58(1H,d,J=7.5Hz), 6.80-7.50(7H,m), 8.05(1H,t,J=1.5Hz), 8.15-8.40(2H,m), 11.56(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )

Table 7 (continued)

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
2-143	165~166	5.10(2H,s), 6.52(1H,d,J=7.5Hz), 7.00-7.21(2H,m), 7.40-7.60(4H,m), 7.72-7.88(2H,m), 8.28(1H,m), 8.68(1H,m), 11.63(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-144	181~182	3.00(6H,s), 6.55(1H,d,J=7.5Hz), 6.73(2H,d,J=9.0Hz), 6.82-7.15(2H,m), 7.20(2H,d,J=9.0Hz), 7.32-7.50(2H,m), 7.74(1H,s), 8.15(2H,m), 11.70(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-145	202~204	6.70-6.85(1H,m), 7.03-7.19(1H,m), 7.38-7.50(1H,m), 7.60-7.83(2H,m), 8.08-8.63(5H,m), 9.67(1H,dd,J=1.35,7.35Hz), 10.49(1H,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
2-146	182~184 (decomp.)	6.35(1H,d,J=7.65Hz), 6.99-7.16(1H,m), 7.21-7.33(1H,m), 7.36-7.66(5H,m), 7.70-7.92(2H,m), 8.58(1H,dd,J=1.35,7.5Hz), 9.83(1H,dd,J=1.35,7.2Hz), 10.15(1H,s), 11.19(1H,d,J=7.65Hz). (CDCl <sub>3</sub> )
2-147	162~164	6.58(1H,d,J=7.5Hz), 6.80-7.50(7H,m), 8.05(1H,d,J=1.5Hz), 8.15-8.40(2H,m), 11.56(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-148	177~179	6.57(1H,d,J=7.8Hz), 6.99-7.12(1H,m), 7.15-7.62(11H,m), 7.68-7.86(2H,m), 8.50(1H,dd,J=1.35,7.5Hz), 8.88(1H,s), 10.30(1H,dd,J=1.35,7.5Hz), 11.42(1H,d,J=7.8Hz). (CDCl <sub>3</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR( $\delta$ value)
2-149	208~210	2.63(3H,s), 6.71(1H,d,J=7.5Hz), 7.00-7.18(1H,m), 7.38-7.50(1H,m), 7.56-7.82(3H,m), 8.14(1H,dd,J=1.35, 8.25Hz), 8.48-8.63(2H,m), 9.58-9.73(1H,m), 10.51(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-150	149~151	2.43(3H,s), 6.60(1H,d,J=7.5Hz), 7.22(1H,t,J=6.75Hz), 7.40-7.80(5H,m), 7.94(1H,s), 8.15(1H,dd,J=6.75, 1.50Hz), 8.85(1H,dd,J=6.75, 1.50Hz), 11.28(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-151	134~135	1.30(3H,t,J=7.0Hz), 2.80(2H,q,J=7.0Hz), 6.40(1H,d,J=7.5Hz), 6.93(1H,t,J=6.75Hz), 7.30-7.75(6H,m), 8.08-8.30(2H,m), 11.52(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-152	141~143	2.45(3H,s), 6.42(1H,d,J=7.5Hz), 7.05-7.38(4H,m), 7.40-7.85(6H,m), 8.13(1H,d,J=6.75Hz), 8.40(1H,s), 8.85(1H,d,J=6.75Hz), 11.18(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-153	235~237	6.40(1H,d,J=7.5Hz), 7.00-7.40(3H,m), 7.45-8.00(7H,m), 8.12(1H,d,J=6.75Hz), 8.50(1H,s), 8.80(1H,d,J=6.75Hz), 11.19(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-154		2.85(3H,s), 6.55(1H,d,J=7.5Hz), 7.09(1H,m), 7.20-7.55(3H,m), 8.60(1H,d,J=6.75Hz), 9.60(1H,d,J=6.75Hz), 10.85(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-155	163~165	6.53(II,d,J=7.5Hz), 6.95(III,t,J=6.75 Hz), 7.00-7.30(II,m), 7.32-7.76(II,m), 7.92(II,s), 8.21(2II,m), 11.57(II,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-156	111~115 (hydrochloride)	6.75(II,d,J=6.75Hz), 6.68-7.11(II,br), 7.11-7.82(7II,m), 7.88-8.11(2II,m), 8.23(II,dd,J=1.35,7.5Hz), 8.68(II,s), 8.92(II,dd,J=1.35,7.5Hz), 11.26(III,d,J=6.75Hz). (DMSO-d <sub>6</sub> )
2-157	174~176	6.52(II,d,J=7.5Hz), 7.00-7.19(2II,m), 7.36-7.58(5II,m), 7.98-8.16(2II,m), 8.25-8.48(2II,m), 11.27(II,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-158	167~168	6.57(II,d,J=8.4Hz), 7.02-7.18(II,m), 7.38-7.70(3II,m), 8.63(II,dd,J=1.35,7.5Hz), 8.66(II,s), 9.61(II,dd,J=1.35,7.5Hz), 10.70(III,d,J=8.4Hz). (CDCl <sub>3</sub> +DMSO-d <sub>6</sub> )
2-159	238~240 (decomp.)	2.58(3II,s), 6.68(III,d,J=6.9Hz), 7.03-7.25(2II,m), 7.46-7.72(3II,m), 8.01-8.24(2II,m), 8.45-8.52(II,m), 8.63(II,s), 8.70-8.86(II,m), 11.11(III,d,J=6.9Hz). (DMSO-d <sub>6</sub> )
2-160	219~220 (decomp.)	6.67(II,d,J=7.2Hz), 6.92-7.10(II,m), 7.27-7.39(II,m), 7.49-7.64(II,m), 7.68-8.02(3II,m), 8.19-8.38(II,m), 8.52(II,dd,J=1.35,7.35Hz), 9.62(II,dd,J=1.35,7.35Hz), 10.37(II,d,J=7.2Hz). (DMSO-d <sub>6</sub> )



Table 7 (continued)

Compound No.	m.p.(°C)	NMR( $\delta$ value)
2-161	265	6.72(1H,d,J=6.6Hz), 7.10-7.57(3H,m), 7.66-(decomp.) 8.01(5H,m), 8.23(1H,dd,J=1.35, 7.5Hz), 8.70(1H,s), 8.91(1H,dd,J=1.35, 7.5Hz), 11.09(1H,d,J=6.6Hz), 13.82(1H,s). (DMSO-d <sub>6</sub> )
2-162	178~179	6.48(1H,d,J=6.5Hz), 7.00(1H,t,J=6.5Hz), 7.00-7.20(1H,m), 8.10(1H,m)8.20-8.35(2H,m), 8.41(1H,s), 11.18(1H,d,J=6.5Hz). (DMSO-d <sub>6</sub> )
2-163	212~213	3.71(3H,s), 3.97(3H,s), 6.58(1H,d,J=6.5Hz), 6.85-7.25(4H,m), 7.45-7.75(3H,m), 8.21(1H,d,J=7.5Hz), 8.40(1H,s), 8.52(1H,d,J=7.5Hz), 11.50(1H,d,J=6.5Hz). (DMSO-d <sub>6</sub> )
2-164	amorphus (triethylamine)	1.36(9H,t,J=7.5Hz), 2.65(3H,s), 3.17(6H,q,J=7.5Hz), 6.56(1H,d,J=7.8Hz), 6.92-7.18(2H,m), 7.28-7.50(2H,m), 8.28(1H,dd,J=1.35, 7.5Hz), 9.00(1H,dd,J=1.35, 7.5Hz), 11.44(1H,d,J=7.8Hz). (CDCl <sub>3</sub> )
2-165	150~152	2.53(3H,s), 2.69(3H,s), 6.55(1H,d,J=7.5Hz), 7.01-7.32(2H,m), 7.36-7.81(3H,m), 8.32-8.50(1H,m), 8.97(1H,dd,J=1.35, 6.6Hz), 11.19(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> +CDCl <sub>3</sub> )
2-166	183~186 (p-toluenesulfonate)	2.29(3H,s), 6.00-6.39(1H,br), 6.73(1H,d,J=7.2Hz), 7.03-7.36(5H,m), 7.40-7.62(6H,m), 7.67-7.78(1H,m), 7.85-8.01(2H,m), 8.19(1H,dd,J=1.35, 7.5Hz), 8.62(1H,s), 8.87(1H,dd,J=1.35, 7.5Hz), 11.19(1H,d,J=7.2Hz). (DMSO-d <sub>6</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR( $\delta$ value)
2-167	195~196	6.73(1H,d,J=7.5Hz), 7.03-7.33(2H,m), 7.36-7.48(1H,m), 7.65(1H,dd,J=1.35, 5.4Hz), 8.17(1H,dd,J=1.35, 7.8Hz), 8.25(1H,s), 8.78(1H,dd,J=1.35, 7.2Hz), 10.42(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-168	253~256	6.55-6.75(3H,m), 7.20(1H,t,J=6.75Hz), 8.00-8.50(5H,m), 8.75-8.90(2H,m), 11.25(1H,d,J=6.0Hz). (DMSO-d <sub>6</sub> )
2-169	196~198	1.15(3H,t,J=7.0Hz), 2.80(3H,s), 3.67(2H,q,J=7.0Hz), 5.74(1H,d,J=9.0Hz), 7.40-7.90(2H,m), 7.60(1H,t,J=6.75Hz), 8.50(1H,d,J=6.75Hz), 9.60(1H,d,J=6.75Hz), 10.50(1H,d,J=9.0Hz). (DMSO-d <sub>6</sub> )
2-170	185~187	2.72(3H,s), 6.52(1H,d,J=7.5Hz), 7.40-7.75(6H,m), 8.50(1H,d,J=6.75Hz), 10.72(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-171	222~223	6.75(1H,d,J=6.5Hz), 7.10-7.30(2H,m), 7.55(1H,m), 7.65-8.80(2H,m), 7.95(1H,d,J=3.5Hz), 8.20(1H,d,J=6.75Hz), 8.64(1H,s), 8.85(1H,d,J=6.75Hz), 10.98(1H,d,J=6.5Hz). (DMSO-d <sub>6</sub> )
2-172	118~120	5.05(2H,br), 6.54(1H,d,J=7.2Hz), 7.21-7.85(6H,m), 8.16-8.31(1H,m), 8.72-8.98(2H,m), 10.48(1H,d,J=7.2Hz). (DMSO-d <sub>6</sub> )
2-173	178~180	6.78(1H,d,J=6.5Hz), 7.20(1H,t,J=6.5Hz), (hydrochloride) 7.30-7.82(7H,m), 8.10(1H,m), 8.56(1H,s), 8.84(1H,m), 11.18(1H,d,J=6.5Hz). (DMSO-d <sub>6</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-174	188~190 (hydrochloride)	6.73(1H,d,J=6.5Hz),6.95(1H,t,J=6.5Hz), 7.00-7.98(8H,m),8.20(1H,m)8.25(1H,s), 8.61(1H,m),11.25(1H,d,J=6.5Hz). (DMSO-d <sub>6</sub> )
2-175	158~159	4.00(3H,s),6.75(1H,d,J=7.5Hz),6.93- 7.45(5H,m),7.55(1H,m),7.75(1H,m), 7.95-8.30(2H,m),8.58(1H,s),8.85(1H,d, J=6.75Hz),11.45(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-176	160~161	3.75(3H,s),6.78(1H,d,J=7.5Hz),6.85- 7.63(7H,m),7.72(1H,dd,J=6.0,1.0Hz), 8.17(1H,dd,J=6.75,1.5Hz),8.59(1H,s), 8.82(1H,dd,6.75,1.5Hz),11.35(1H,d,J= 7.5Hz).(DMSO-d <sub>6</sub> )
2-177	229~230	3.86(3H,s),6.74(1H,d,J=7.5Hz),6.90- 7.25(3H,m),7.50(1H,m),7.72(1H,dd,J= 6.0,1.0Hz),7.88(2H,d,J=8.5Hz),8.12 (1H,dd,J=6.75,1.5Hz),8.42(1H,s),8.88 (1H,dd,J=6.75,1.5Hz),11.40(1H,d,J= 7.5Hz).(DMSO-d <sub>6</sub> )
2-178	172~173	6.78(1H,d,J=7.5Hz),7.15-7.85(6H,m), 8.00-8.35(3H,m),8.85(1H,s),8.95(1H,d, J=6.75Hz),11.25(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-179	225~226	6.75(1H,d,J=7.5Hz),7.05-7.30(2H,m), 7.40-7.60(3H,m),7.65-8.05(3H,m),8.18 (1H,d,J=6.75Hz),8.65(1H,s),8.81(1H,d, J=6.75Hz),11.23(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-180	228~229	6.75(1H,d,J=7.5Hz), 7.00-7.30(2H,m), 7.40-7.65(2H,m), 7.77(1H,dd,J=6.0, 1.0Hz), 7.95(2H,d,J=8.5Hz), 8.15(1H,dd, J=6.75,1.5Hz), 8.63(1H,s), 8.84(1H,dd, J=6.75,1.5Hz), 11.35(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-181	179~180	2.37(3H,s), 2.48(3H,s), 6.67(1H,d,J= 7.5Hz), 6.91-7.18(4H,m), 7.33-7.45(2H, m), 7.60(1H,d,J=9.0Hz), 7.78(1H,s), 8.28-8.40(2H,m), 11.53(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-182	193~195	2.30(6H,s), 6.50(1H,d,J=7.5Hz), 6.91 (1H,t,J=6.0Hz), 7.03-7.30(3H,m), 7.39- 7.51(2H,m), 7.53-7.65(1H,m), 7.85(1H,s) , 8.10-8.28(2H,m), 11.60(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-183	205~207	2.50(6H,br), 6.75(1H,d,J=8.5Hz), 7.00- 7.30(4H,m), 7.38-7.53(1H,m), 7.65-7.85 (2H,m), 8.14(1H,d,J=6.75Hz), 8.35(1H,s) , 8.85(1H,d,J=6.75Hz), 11.28(1H,d,J= 8.5Hz). (DMSO-d <sub>6</sub> )
2-184	103~104	1.34(9H,s), 2.50(3H,s), 6.60(1H,d,J= 8.5Hz), 6.80-7.15(2H,m), 7.20-7.45(4H, m), 7.60-7.85(2H,m), 8.18(1H,d,J= 6.75Hz), 8.30(1H,d,J=6.75Hz), 11.59(1H, d,J=8.5Hz). (CDCl <sub>3</sub> )
2-185	129~130	1.32(9H,s), 2.45(3H,s), 6.65(1H,d,J= 8.5Hz), 6.80-7.20(2H,m), 7.21-7.51(4H, m), 7.75-7.90(2H,m), 8.21(1H,d,J= 6.75Hz), 8.32(1H,d,J=6.75Hz), 11.55(1H, d,J=8.5Hz). (CDCl <sub>3</sub> )

Table 7 (continued)

Compound No.	m.p. (°C)	NMR ( $\delta$ value)
2-186	159~160	2.35(3H,s), 6.75(1H,d,J=8.5Hz), 7.08-7.30(3H,m), 7.38-7.56(2H,m), 7.72(1H,d,J=4.5Hz), 7.90(1H,m), 8.18(1H,d,J=6.75Hz), 8.78(1H,s), 8.92(1H,d,J=6.75Hz), 11.18(1H,d,J=8.5Hz). (DMSO-d <sub>6</sub> )
2-187	184~185	2.40(3H,s), 6.75(1H,d,J=7.5Hz), 7.00-7.90(8H,m), 8.18(1H,d,J=6.75Hz), 8.58(1H,s), 8.85(1H,d,J=6.75Hz), 11.43(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-188	218~220	2.35(3H,s), 6.76(1H,d,J=7.5Hz), 7.00-7.40(4H,m), 7.54(1H,m), 7.65-7.95(3H,m), 8.13(1H,dd,J=6.75,1.5Hz), 8.50(1H,s), 8.83(1H,dd,J=6.75,1.5Hz), 11.38(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-189	186~188 (decomp.)	2.67(3H,s), 6.61(1H,d,J=7.8Hz), 6.92-7.16(2H,m), 7.33-7.67(6H,m), 7.86-8.07(2H,m), 8.23(1H,dd,J=1.35,6.45Hz), 8.61(1H,dd,J=1.35,6.45Hz), 8.71(1H,s), 11.33(1H,d,J=7.8Hz). (CDCl <sub>3</sub> )
2-190	132~134	6.49(1H,d,J=7.8Hz), 6.90-7.50(10H,m), 8.10-8.30(2H,m), 10.71(1H,d,J=7.8Hz). (CDCl <sub>3</sub> )
2-191	155~157	6.71(1H,d,J=7.5Hz), 6.99-7.19(2H,m), 7.20-7.53(6H,m), 7.60(1H,dd,J=1.35,5.1Hz), 8.13(1H,dd,J=1.35,7.5Hz), 8.38(1H,s), 8.88(1H,dd,J=1.35,6.75Hz), 10.80(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR( $\delta$ value)
2-194	235~237	6.75(1H,d,J=6.3Hz), 7.05-7.34(2H,m), 7.50-7.60(3H,m), 7.78-8.07(5H,m), 8.15 (1H,dd,J=1.5,6.45Hz), 8.39-8.47(1H,m), 8.69(1H,s), 8.83(1H,dd,J=1.5, 6.45Hz), 11.38(1H,d,J=6.3Hz). (DMSO-d <sub>6</sub> )
2-195	170~172	2.37(3H,s), 3.02(2H,br), 6.60(1H,d,J= 8.1Hz), 6.70-7.14(2H,m), 7.35-7.51(2H, m), 8.00-8.29(2H,m), 11.49(1H,d,J= 8.1Hz). (CDCl <sub>3</sub> )
2-196	187~189	2.20-2.45(12H,m), 6.74(1H,d,J=8.5Hz), 7.00-7.28(2H,m), 7.35(1H,s), 7.45(1H,d, J=3.0Hz), 7.63(1H,d,J=6.0Hz), 8.05- 8.28(2H,m), 8.82(1H,d,J=6.75Hz), 11.40 (1H,d,J=8.5Hz). (DMSO-d <sub>6</sub> )
2-197	118~120	1.35(3H,t,J=7.0Hz), 2.76(2H,q,J=7.0Hz) , 6.62(1H,d,J=7.5Hz), 6.85-7.10(2H,m), 7.35(1H,m), 7.65(1H,d,J=4.5Hz), 7.75(1 H,s), 8.98(1H,d,J=6.75, 1.5Hz), 8.62(1H, d,J=6.75Hz), 11.27(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-200	195~196	1.23(3H,t,J=7.0Hz), 2.66(2H,q,J=7.0Hz) , 6.71(1H,d,J=8.5Hz), 7.00-7.40(4H,m), 7.52(1H,d,J=3.0Hz), 7.60-7.90(3H,m), 8.10(1H,d,J=6.75Hz), 8.43(1H,s), 8.75 (1H,d,J=6.75Hz), 11.40(1H,d,J=8.5Hz). (DMSO-d <sub>6</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-201	237~238	6.72(1H,d,J=8.5Hz), 7.05-7.27(2H,m), 7.45-7.95(6H,m), 8.13(1H,dd,J=6.75, 1.5Hz), 8.55(1H,s), 8.80(1H,dd,J=6.75, 1.5Hz), 11.26(1H,d,J=8.5Hz). (DMSO-d <sub>6</sub> )
2-202	91~94	1.03-1.37(6H,m), 1.47-3.00(4H,m), 6.82(1H,d,J=7.5Hz), 7.05-7.31(4H,m), 7.38-7.49(1H,m), 7.52-7.61(1H,m), 7.63-7.76(1H,m), 8.15(1H,d,J=1.35, 7.5Hz), 8.38(1H,s), 8.88(1H,dd,J=1.35, 6.95Hz), 11.21(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> )
2-203	189~190	2.54(3H,s), 6.75(1H,d,J=8.5Hz), 7.05-7.40(4H,m), 7.48(1H,m), 7.70(1H,m), 7.93(1H,m), 8.25(1H,d,J=6.75Hz), 8.50(1H,s), 8.85(1H,d,J=6.75Hz), 11.12(1H,d,J=8.5Hz). (DMSO-d <sub>6</sub> )
2-204	141~142	1.18(3H,t,J=7.8Hz), 2.44(3H,s), 2.60(2H,q,J=7.8Hz), 6.77(1H,d,J=7.2Hz), 7.02-7.30(3H,m), 7.40-7.52(1H,m), 7.63-7.77(2H,m), 8.13(1H,dd,1.35, 7.5Hz), 8.41(1H,s), 8.83(1H,dd,J=1.35, 6.75Hz), 11.28(1H,d,J=7.2Hz). (DMSO-d <sub>6</sub> )
2-209	195~197	2.62(3H,s), 2.81(3H,s), 6.65(1H,d,J=8.5Hz), 6.90-7.50(5H,m), 7.80-8.15(3H,m), 8.42(1H,s), 8.78(1H,d,J=6.75Hz), 11.08(1H,d,J=8.5Hz). (DMSO-d <sub>6</sub> )

Table 7 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-210	237~238	2.61(3H,s), 3.22(3H,s), 6.74(1H,d,J=8.5Hz), 7.05-7.35(2H,m), 7.48(1H,m), 7.65-8.22(5H,m), 8.58(1H,s), 8.88(1H,d,J=6.75Hz), 11.18(1H,d,J=8.5Hz). (DMSO-d <sub>6</sub> )
2-211	201~202	2.30(3H,s), 3.94(3H,s), 6.49(1H,d,J=7.5Hz), 6.91(1H,t,J=7.5Hz), 7.00-7.30(3H,m), 7.35-7.52(3H,m), 7.89(1H,d,J=2.0Hz), 8.22(1H,s), 8.18(1H,dd,J=7.5, 1.5Hz), 8.25(1H,dd,J=7.5, 1.5Hz), 11.62(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-212	244~245	4.00(3H,s), 6.50(1H,d,J=7.5Hz), 6.99-7.35(3H,m), 7.48-7.61(2H,m), 8.05-8.25(2H,m), 8.42(1H,s), 8.59(1H,dd,J=7.5, 1.5Hz), 11.29(1H,d,J=7.5Hz). (DMSO-d <sub>6</sub> +CDCl <sub>3</sub> )
2-213	138~139	1.10-1.40(12H,m), 2.75-3.15(1H,m), 3.55-3.90(1H,m), 6.82(1H,d,J=8.5Hz), 7.05-7.70(7H,m), 8.18(1H,d,J=6.75Hz), 8.23(1H,s), 8.89(1H,d,J=6.75Hz), 11.20(1H,d,J=8.5Hz). (DMSO-d <sub>6</sub> )
2-214	153~154	1.12-1.31(6H,m), 1.30(3H,d,J=6.0Hz), 2.70-3.18(1H,m), 3.40-3.70(1H,m), 6.67(1H,d,J=8.5Hz), 6.90-7.20(2H,m), 7.70(1H,s), 8.15-8.40(2H,m), 11.42(1H,d,J=8.5Hz). (CDCl <sub>3</sub> )



Table 7 (continued)

Compound No.	m.p.(°C)	NMR ( $\delta$ value)
2-215	207~209	1.32(3H,d,J=7.5Hz), 2.75-3.15(1H,m), 6.55(1H,d,J=7.5Hz), 6.98(1H,t,J=6.0Hz), 7.08-7.18(1H,m), 7.28(1H,d,J=8.5Hz), 7.35-7.55(2H,m), 7.75(2H,d,J=8.5Hz), 7.89(1H,s), 8.20(1H,dd,J=6.0, 1.5Hz), 8.25(1H,dd,J=6.0, 1.5Hz), 11.60(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-231	175~176	6.50(1H,d,J=7.5Hz), 7.05-7.15(1H,m), 7.35-7.50(4H,m), 7.80-8.00(1H,m), 7.85(1H,d,J=4.5Hz), 8.32(1H,s), 8.50(1H,d,J=4.5Hz), 11.12(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )
2-232	209~210	6.48(1H,d,J=7.5Hz), 7.05-7.20(1H,m), 7.35-7.55(4H,m), 7.75-7.95(1H,m), 7.82(1H,s), 8.38(1H,s), 10.98(1H,d,J=7.5Hz). (CDCl <sub>3</sub> )

## Formulation Example 1

Emulsifiable concentrates, as prepared by blending the following components:

Compound No. 2-2	20% (by weight)
xylene	75%
polyoxyethylenealkylaryl ether (Nonipol 85 <sup>R</sup> )	5%

## Formulation Example 2

Wettable powders, as prepared by blending and grinding the following components:

Compound No. 2-24	50% (by weight)
diatomaceous earth	44%
polyoxyethylenealkylaryl ether (Nonipol 85 <sup>R</sup> )	6%

## Formulation Example 3

Powders, as prepared by blending the following mixture:

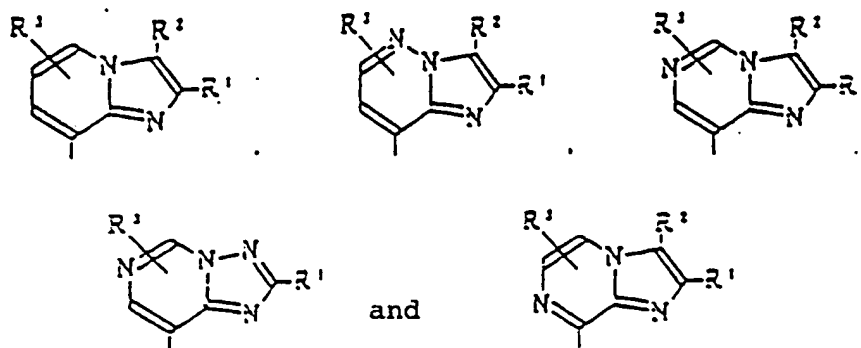
Compound No. 2-27	3% (by weight)
clay	40%
talc	57%

## Claims

Claims for the following Contracting States : AT, BE, CH, DE, DK, FR, GB, GR, IT, LI, LU, NL, SE

1. A condensed heterocyclic compound of the formula: or an agriculturally acceptable salt thereof, wherein:

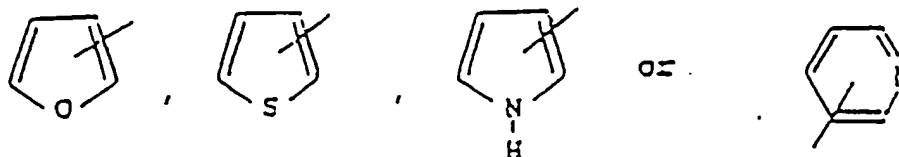
(1) Q is a condensed heterocyclic group having a nitrogen atom in the bridgehead selected from the group consisting of



in which R<sup>1</sup> is C<sub>1-6</sub> alkyl, halogen, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio, C<sub>6-10</sub> aryloxy, C<sub>6-10</sub> arylthio, alkoxycarbonyl, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group;

R<sup>2</sup> and R<sup>3</sup> are hydrogen, C<sub>1-6</sub> alkyl, halogen, nitro, amino, sulfo, mono- or dialkylsulfamoyl, alkoxycarbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group;

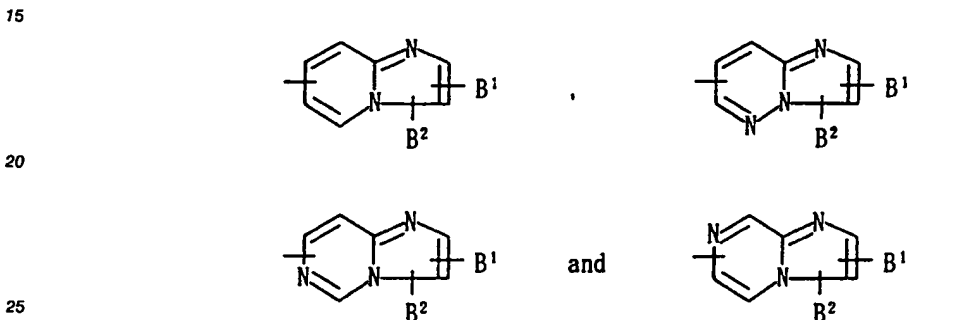
X is hydrogen; a group attached through a carbon atom selected from the group consisting of C<sub>1-10</sub> alkyl, C<sub>2-4</sub> alkenyl, C<sub>1-10</sub> haloalkyl, C<sub>3-6</sub> cycloalkyl, C<sub>3-6</sub> cycloalkenyl, C<sub>6-10</sub> aryl, C<sub>7-10</sub> aralkyl and a heterocyclic ring of the formula



a group attached through an oxygen atom selected from the group consisting of C<sub>1-10</sub> alkoxy, C<sub>6-10</sub> aryloxy and C<sub>7-10</sub> aralkyloxy; a group attached through a sulfur atom selected from the group consisting of C<sub>1-10</sub> alkylthio, C<sub>6-10</sub> arylthio and C<sub>7-10</sub> aralkylthio; a group attached through a nitrogen atom selected from the group consisting of C<sub>1-10</sub> alkylamino, C<sub>6-10</sub> arylamino, C<sub>7-10</sub> aralkylamino and a heterocyclic ring of the formula



10 wherein each of said groups attached through a carbon atom, an oxygen atom, a sulfur atom or a nitrogen atom is unsubstituted or substituted with one to four substituents selected from the group consisting of nitro, amino, hydroxy, cyano, carboxyl, sulfonyl, halogen, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio, phenylthio and benzylthio; and Y is cyano, carbamoyl, thiocarbamoyl or trichloromethyl group; or (2) Q is a condensed heterocyclic group having a nitrogen atom in the bridgehead selected from the group consisting of



30 in which the substituents B<sup>1</sup> and B<sup>2</sup> on the condensed heterocyclic group for Q may be the same or different each other, and are hydrogen atom, nitro, amino, hydroxyl, cyano, C<sub>1-3</sub> acyl, carbamoyl, carboxyl, C<sub>1-4</sub> alkoxycarbonyl, sulfo, halogen, C<sub>1-4</sub> alkoxy, C<sub>6-10</sub> aryloxy, C<sub>6-10</sub> arylcarbonyl, C<sub>6-10</sub> aryl, C<sub>7-10</sub> aralkyl, C<sub>3-7</sub> cycloalkyl, C<sub>1-4</sub> alkylthio, C<sub>6-10</sub> arylthio, C<sub>7-10</sub> aralkylthio, mono- or di-C<sub>1-4</sub> alkylsulfamoyl, C<sub>1-6</sub> alkyl, mono-, di-, tri- or tetra-halogen-substituted phenyl, mono-, di-, tri- or tetra-C<sub>1-4</sub> alkyl-substituted phenyl, mono-, di-, tri- or tetra-C<sub>1-4</sub> alkoxy-substituted phenyl, mono-, di-, tri- or tetra-C<sub>1-4</sub> alkylthio-substituted phenyl, 2-chloro-4-nitrophenyl, 4-nitrophenyl, 2-methyl-4-aminophenyl, 2-bromo-4-nitrophenyl, 2-nitro-4-methylphenyl, or 5 or 6 membered heterocyclic group selected from pyridyl, furyl, thienyl and thiazolyl; and X and Y have the same meaning as defined above; or (3) Q is:

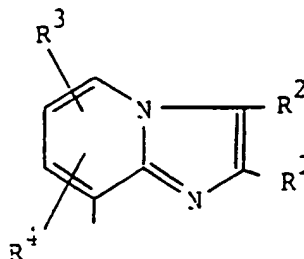


and

50 R<sup>1</sup> is methyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is 1,1,2,2-tetrafluoroethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is morpholinomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is isopropylthiomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is 2-cyanoethyl, X is 3-fluorophenyl and Y is cyano;  
 55 or R<sup>1</sup> is methyl, R<sup>2</sup> is 2-methyl-1-propenyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 2-N-methylpyrrolyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 3-trifluoromethylphenyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 3-methylphenyl and Y is cyano;

or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 5-methylfur-2-yl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 2-pyridylmethyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 3-methylthien-2-yl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 5 or R<sup>1</sup> is phenyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is 1,1,2,2-tetrafluoroethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is 2-cyanoethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is 2-methyl-1-propenyl, X is 2-furyl and Y is cyano;  
 10 or R<sup>1</sup> is phenyl, R<sup>2</sup> is 2,2,2-trichloro-1-hydroxyethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is trifluoromethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-1-propenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is ethoxyiminomethyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenoxyethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 15 or R<sup>1</sup> is phenylthiomethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenylsulfinylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenylsulfonylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenoxyethyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is methoxymethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 20 or R<sup>1</sup> is methoxymethyl, R<sup>2</sup> is morpholinomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylthiomethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylsulfinylmethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylthiomethyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylsulfonylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 25 or R<sup>1</sup> is morpholinomethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is diethoxyphosphorylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is (dimethylthiocarbonyl)thiomethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is chloromethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is trifluoromethyl, R<sup>2</sup> is phenylcarbamoyl, X is 3-fluorophenyl and Y is cyano;  
 30 or R<sup>1</sup> is hydrogen, R<sup>2</sup> is 1-hydroxybenzyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is sulfomethyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 4-dimethylaminophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is benzylidenamino, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is biphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 35 or R<sup>1</sup> is 4-sulfophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is trifluoromethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is formamide;  
 or R<sup>1</sup> is 3-hydroxyphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-hydroxyphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is benzylidenamino, X is 2-thienyl and Y is cyano;  
 40 or R<sup>1</sup> is 2-dimethylcarbamoylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methoxymethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-4-methylsulfinylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-4-methylsulfonylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 4-cyanophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 45 or R<sup>1</sup> is 2-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 3-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-5-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;

or (4) Q is:



and

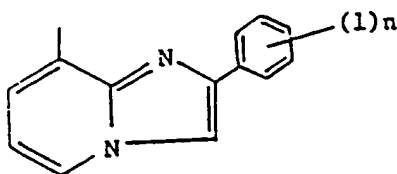
R¹ is methyl, R² is N,N-dimethylaminomethyl, R³ is 5-methyl, R⁴ is hydrogen, X is 2-thienyl and Y is cyano;

or R¹ is phenyl, R² is hydrogen, R³ is 5-methyl, R⁴ is 7-methyl, X is 2-furyl and Y is cyano;

or R¹ is phenyl, R² is hydrogen, R³ is 5-trifluoromethyl, R⁴ is 7-trifluoromethyl, X is 2-furyl and Y is cyano;

or R¹ is phenyl, R² is hydrogen, R³ is 5-hydroxy, R⁴ is 6-chloro, X is 2-thienyl and Y is cyano.

2. A compound or claim 1, wherein Q is



wherein n is 0, 1, 2, 3 or 4 and 1 is the same or different and selected from a group comprising a C₁-₄ alkyl, C₁-₄ alkoxy, C₁-₄ alkylthio or halogen,

X is phenyl, fluorine-substituted phenyl, thienyl or furyl group, and

Y is cyano,

or its salt.

3. A compound of claim 1, which is α-(2-phenylimidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)acetonitrile or its salt.

4. A compound of claim 1, which is α-{2-(2-methoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.

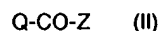
5. A compound of claim 1, which is α-{2-(3-methoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.

6. A compound of claim 1, which is α-{2-(2,4-dimethoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.

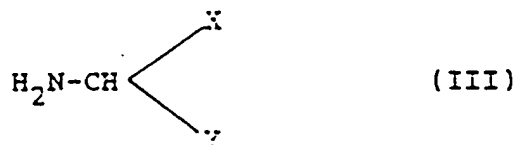
7. A compound of claim 1, which is α-{2-(2,4-diisopropylphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.

8. A compound of claim 1, which is α-{2-(3-trifluoromethylphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.

9. A method for preparing a condensed heterocyclic compound of the formula (I) in claim 1 or its salt which comprises reacting a compound of the formula



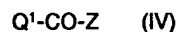
wherein Q has the same meaning as defined in claim 1 and COZ denotes a carboxylic group or its reactive derivative or its salt with a compound of the general formula



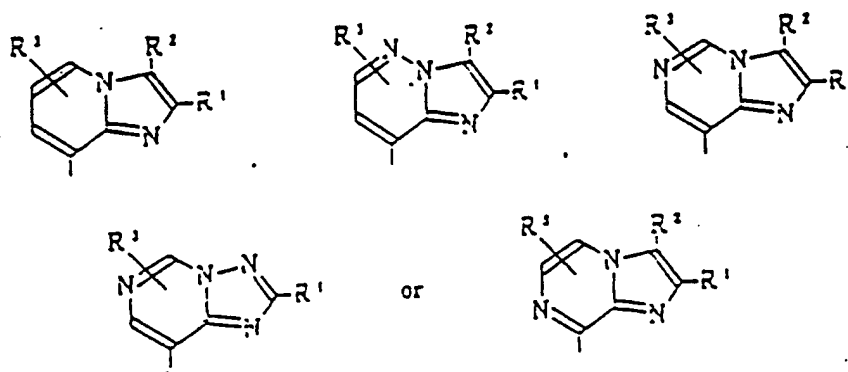
wherein X and Y are each as defined in claim 1, or its salt.

10. A fungicidal composition comprising a condensed heterocyclic compound of the formula (I) in claim 1 or its salt, and a carrier or diluent.

11. A condensed heterocyclic carboxylic acid compound of the formula:



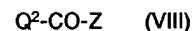
wherein Q<sup>1</sup> is a condensed heterocyclic group having a nitrogen atom in the bridgehead as follows



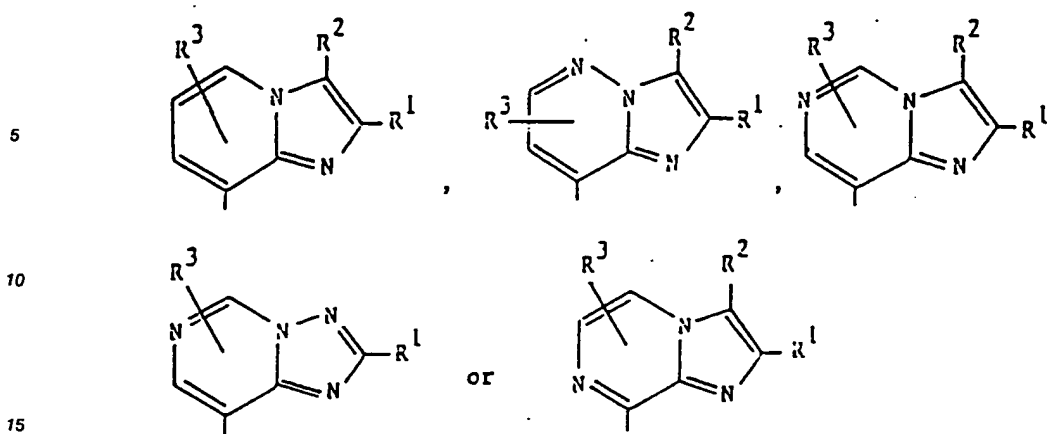
in which R<sup>1</sup> is an C<sub>1-6</sub> alkyl, halogen, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio, C<sub>6-10</sub> aryloxy, C<sub>6-10</sub> arylthio, alkoxycarbonyl, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group;

R<sup>2</sup> and R<sup>3</sup> are hydrogen, C<sub>1-6</sub> alkyl, halogen, nitro, amino, sulfo, mono- or dialkylsulfamoyl, alkoxycarbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group, and COZ denotes a carboxylic group or its reactive derivative; or its salt.

12. A process for preparing a compound of the formula



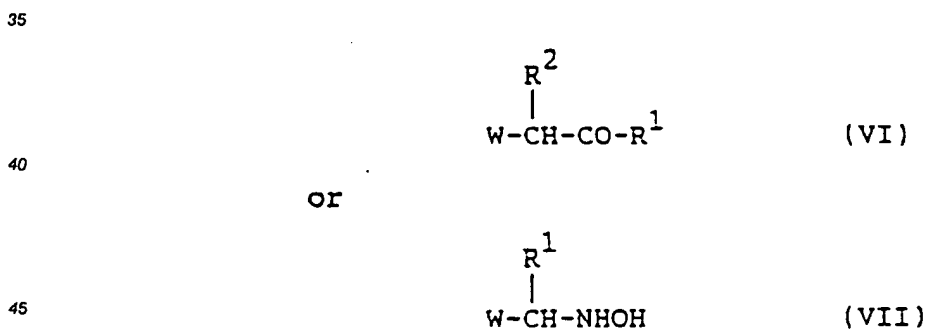
wherein Q<sup>2</sup> is a group of the general formula:



in which R<sup>1</sup>, R<sup>2</sup> and R<sup>3</sup> have the same meanings as defined in claim 11 or its salt which comprises reacting a compound of the formula:



wherein the ring  $\textcircled{A}$  is a pyridine, pyridazine, pyrimidine or pyrazine group which is unsubstituted or substituted by an C<sub>1-6</sub> alkyl, halogen, nitro, amino, sulfo, mono- or di-alkylsulfamoyl, alkoxy, carbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group, and COZ denotes a carboxylic group or its reactive derivative, or its salt with a compound of the formula:

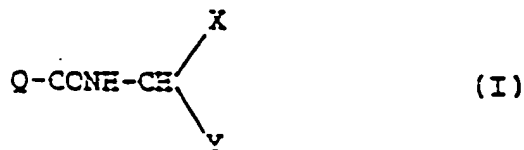


wherein W is a halogen atom, and R<sup>1</sup> and R<sup>2</sup> have the same meanings as defined above.

- 50 13. Use of a compound as claimed in claims 1 to 8 as fungicide more particularly for controlling various kinds of plant blight such as downy mildew and blight.

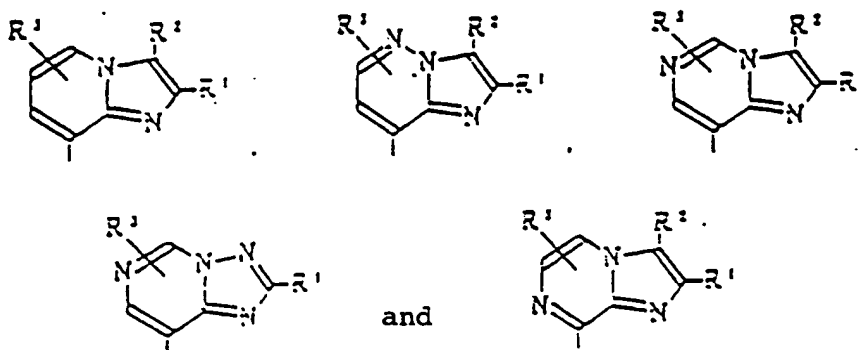
## Claims for the following Contracting State : ES

1. A method for preparing a condensed heterocyclic compound of the formula (I):



or an agriculturally acceptable salt thereof, wherein:

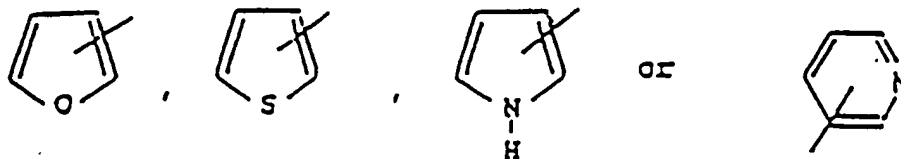
- (1) Q is a condensed heterocyclic group having a nitrogen atom in the bridgehead selected from the group consisting of



in which R<sup>1</sup> is C<sub>1-6</sub> alkyl, halogen, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio, C<sub>6-10</sub> arylloxy, C<sub>6-10</sub> arylthio, alkoxycarbonyl, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group;

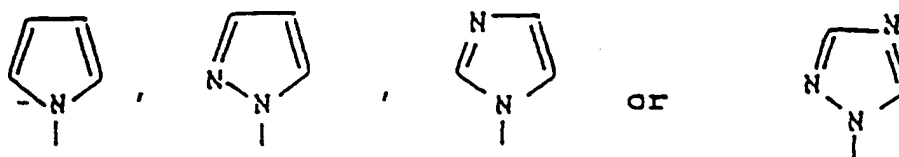
R<sup>2</sup> and R<sup>3</sup> are hydrogen, C<sub>1-6</sub> alkyl, halogen, nitro, amino, sulfo, mono- or dialkylsulfamoyl, alkoxycarbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group;

X is hydrogen; a group attached through a carbon atom selected from the group consisting of C<sub>1-10</sub> alkyl, C<sub>2-4</sub> alkenyl, C<sub>1-10</sub> haloalkyl, C<sub>3-6</sub> cycloalkyl, C<sub>3-6</sub> cycloalkenyl, C<sub>6-10</sub> aryl, C<sub>7-10</sub> aralkyl and a heterocyclic ring of the formula

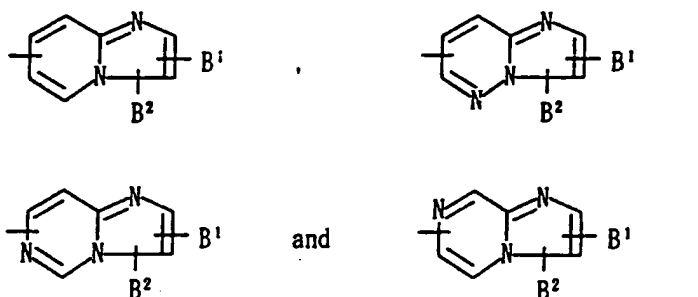


a group attached through an oxygen atom selected from the group consisting of C<sub>1-10</sub> alkoxy, C<sub>6-10</sub> aryloxy and C<sub>7-10</sub> aralkyloxy; a group attached through a sulfur atom selected from the group consisting of C<sub>1-10</sub> alkylthio, C<sub>6-10</sub> arylthio and C<sub>7-10</sub> aralkylthio; a group attached through a nitrogen atom selected from the group consisting of C<sub>1-10</sub> alkylamino, C<sub>6-10</sub> arylamino, C<sub>7-10</sub> aralkylamino and a heterocyclic ring of the formula





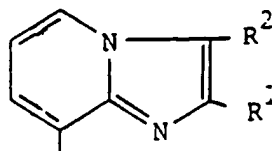
10 wherein each of said groups attached through a carbon atom, an oxygen atom, a sulfur atom or a nitrogen atom is unsubstituted or substituted with one to four substituents selected from the group consisting of nitro, amino, hydroxy, cyano, carboxyl, sulfonyl, halogen, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio, phenylthio and benzylthio; and Y is cyano, carbamoyl, thiocarbamoyl or trichloromethyl group; or (2) Q is a condensed heterocyclic group having a nitrogen atom in the bridgehead selected from the group consisting of



30 in which the substituents B<sup>1</sup> and B<sup>2</sup> on the condensed heterocyclic group for Q may be the same or different each other, and are hydrogen atom, nitro, amino, hydroxyl, cyano, C<sub>1-3</sub> acyl, carbamoyl, carboxyl, C<sub>1-4</sub> alkoxy, C<sub>6-10</sub> aryloxy, C<sub>6-10</sub> arylcarbonyl, C<sub>6-10</sub> aryl, C<sub>7-10</sub> aralkyl, C<sub>3-7</sub> cycloalkyl, C<sub>1-4</sub> alkylthio, C<sub>6-10</sub> arylthio, C<sub>7-10</sub> aralkylthio, mono- or di-C<sub>1-4</sub> alkylsulfamoyl, C<sub>1-6</sub> alkyl, mono-, di-, tri- or tetra-halogen-substituted phenyl, mono-, di-, tri- or tetra-C<sub>1-4</sub> alkyl-substituted phenyl, mono-, di-, tri- or tetra-C<sub>1-4</sub> alkoxy-substituted phenyl, mono-, di-, tri- or tetra-C<sub>1-4</sub> alkylthio-substituted phenyl, 2-chloro-4-nitrophenyl, 4-nitrophenyl, 2-methyl-4-aminophenyl, 2-bromo-4-nitrophenyl, 2-nitro-4-methylphenyl, or 5 or 6 membered heterocyclic group selected from pyridyl, furyl, thienyl and thiazolyl; and X and Y have the same meaning as defined above; or (3) Q is:

35

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and

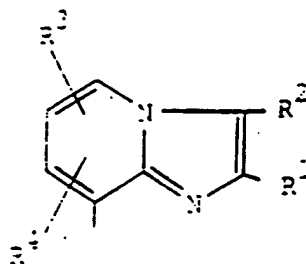
50 R<sup>1</sup> is methyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is 1,1,2,2-tetrafluoroethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is morpholinomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is isopropylthiomethyl, X is 3-fluorophenyl and Y is cyano;  
 55 or R<sup>1</sup> is methyl, R<sup>2</sup> is 2-cyanoethyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is methyl, R<sup>2</sup> is 2-methyl-1-propenyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 2-N-methylpyrrolyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 3-trifluoromethylphenyl and Y is cyano;

or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 3-methylphenyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 5-methylfur-2-yl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 2-pyridylmethyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, X is 3-methylthien-2-yl and Y is cyano;  
 5 or R<sup>1</sup> is phenyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is 1,1,2,2-tetrafluoroethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is 2-cyanoethyl, X is 2-furyl and Y is cyano;  
 10 or R<sup>1</sup> is phenyl, R<sup>2</sup> is 2-methyl-1-propenyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is 2,2,2-trichloro-1-hydroxyethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is trifluoromethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-1-propenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is ethoxyiminomethyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 15 or R<sup>1</sup> is phenoxyimethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenylthiomethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenylsulfinylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenylsulfonylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is phenoxyimethyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 3-fluorophenyl and Y is cyano;  
 20 or R<sup>1</sup> is methoxymethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methoxymethyl, R<sup>2</sup> is morpholinomethyl, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylthiomethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylsulfinylmethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is methylthiomethyl, R<sup>2</sup> is N,N-dimethylaminomethyl, X is 2-furyl and Y is cyano;  
 25 or R<sup>1</sup> is methylsulfonylmethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is morpholinomethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is diethoxyphosphoxymethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is (dimethylthiocarbamoyl)thiomethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 or R<sup>1</sup> is chloromethyl, R<sup>2</sup> is hydrogen, X is 2-furyl and Y is cyano;  
 30 or R<sup>1</sup> is trifluoromethyl, R<sup>2</sup> is phenylcarbamoyl, X is 3-fluorophenyl and Y is cyano;  
 or R<sup>1</sup> is hydrogen, R<sup>2</sup> is 1-hydroxybenzyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is sulfomethyl, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 4-dimethylaminophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is phenyl, R<sup>2</sup> is benzylidenamino, X is 2-thienyl and Y is cyano;  
 35 or R<sup>1</sup> is biphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 4-sulfophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is trifluoromethyl, R<sup>2</sup> is hydrogen, X is 3-fluorophenyl and Y is formamide;  
 or R<sup>1</sup> is 3-hydroxyphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-hydroxyphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 40 or R<sup>1</sup> is methyl, R<sup>2</sup> is benzylidenamino, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-dimethylcarbamoylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methoxymethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-4-methylsulfinylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-4-methylsulfonylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 45 or R<sup>1</sup> is 4-cyanophenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 3-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;  
 or R<sup>1</sup> is 2-methyl-5-trifluoromethylphenyl, R<sup>2</sup> is hydrogen, X is 2-thienyl and Y is cyano;

50

55

or (4) Q is:



and

R<sup>1</sup> is methyl, R<sup>2</sup> is N,N-dimethylaminomethyl, R<sup>3</sup> is 5-methyl, R<sup>4</sup> is hydrogen, X is 2-thienyl and Y is cyano;

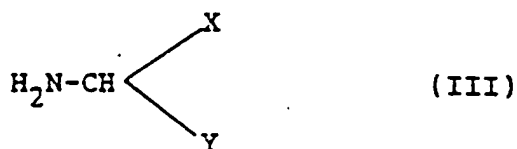
or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, R<sup>3</sup> is 5-methyl, R<sup>4</sup> is 7-methyl, X is 2-furyl and Y is cyano;

or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, R<sup>3</sup> is 5-trifluoromethyl, R<sup>4</sup> is 7-trifluoromethyl, X is 2-furyl and Y is cyano;

or R<sup>1</sup> is phenyl, R<sup>2</sup> is hydrogen, R<sup>3</sup> is 5-hydroxy, R<sup>4</sup> is 6-chloro, X is 2-thienyl and Y is cyano which comprises reacting a compound of the formula

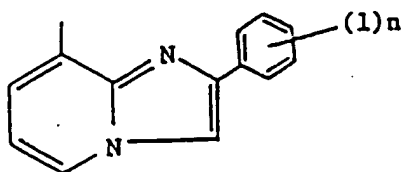
Q-CO-Z (II)

wherein Q has the same meaning as defined above and COZ denotes a carboxylic group or its reactive derivative or its salt with a compound of the general formula



wherein X and Y are each as defined above, or its salt.

2. A method of claim 1, wherein Q is



wherein n is 0, 1, 2, 3 or 4 and L is the same or different and selected from a group comprising a C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio or halogen,

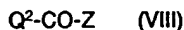
X is phenyl, fluorine-substituted phenyl, thienyl or furyl group, and

Y is cyano, or its salt.

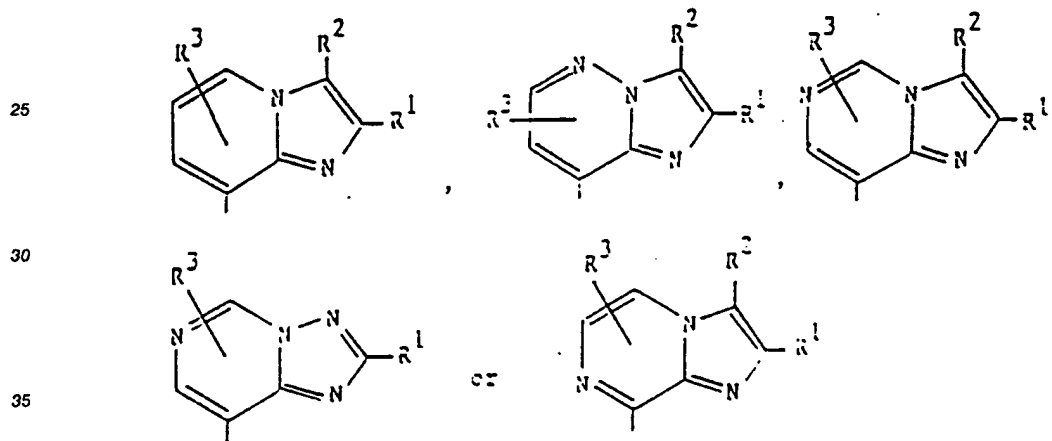
3. A method for preparing a compound according to claim 1, which is  $\alpha$ -(2-phenylimidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)acetonitrile or its salt.

4. A method for preparing a compound according to claim 1, which is  $\alpha$ -[2-(2-methoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino]-(2-thienyl)acetonitrile or its salt.

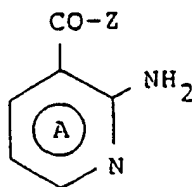
5. A method for preparing a compound according to claim 1, which is  $\alpha$ -{2-(3-methoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.
6. A method for preparing a compound according to claim 1, which is  $\alpha$ -{2-(2,4-dimethoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.
7. A method for preparing a compound according to claim 1, which is  $\alpha$ -{2-(2,4-diisopropylphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.
8. A method for preparing a compound according to claim 1, which is  $\alpha$ -{2-(3-trifluoromethylphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitrile or its salt.
9. A method for preparing a fungicidal composition comprising mixing a condensed heterocyclic compound of the formula (I) in claim 1 or its salt, and a carrier or diluent.
10. A process for preparing a compound of the formula



wherein  $Q^2$  is a condensed heterocyclic group having a nitrogen atom in the bridgehead as follows

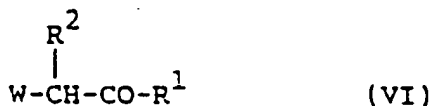


in which  $R^1$  is an  $C_{1-6}$  alkyl, halogen,  $C_{1-4}$  alkoxy,  $C_{1-4}$  alkylthio,  $C_{6-10}$  aryloxy,  $C_{6-10}$  arylthio, alkoxycarbonyl, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen,  $C_{1-4}$  alkyl,  $C_{1-4}$  alkoxy,  $C_{1-4}$  alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group;  $R^2$  and  $R^3$  are a hydrogen,  $C_{1-6}$  alkyl, halogen, nitro, amino, sulfo, mono- or di-alkylsulfamoyl, alkoxycarbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen,  $C_{1-4}$  alkyl,  $C_{1-4}$  alkoxy,  $C_{1-4}$  alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group, and COZ denotes a carboxylic group or its reactive derivative; or its salt which comprises reacting a compound of the formula:

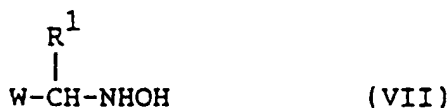


wherein the ring  $\textcircled{A}$  is a pyridine, pyridazine, pyrimidine or pyrazine group which is unsubstituted or

substituted by an C<sub>1-6</sub> alkyl, halogen, nitro, amino, sulfo, mono- or di-alkylsulfamoyl, alkoxycarbonyl, formyl, cyano, phenyl, phenyl substituted with 1 to 4 group(s), which may be the same or different and are selected from the group consisting of halogen, C<sub>1-4</sub> alkyl, C<sub>1-4</sub> alkoxy, C<sub>1-4</sub> alkylthio and nitro or 5 or 6 membered aromatic heterocyclic group, and COZ denotes a carboxylic group or its reactive derivative, or its salt with a compound of the formula:



or



wherein W is a halogen atom, and R<sup>1</sup> and R<sup>2</sup> have the same meanings as defined above.

11. Use of a compound as prepared according to claims 1 to 8 as fungicide more particularly for controlling various kinds of plant blight such as downy mildew and blight.

#### Patentansprüche

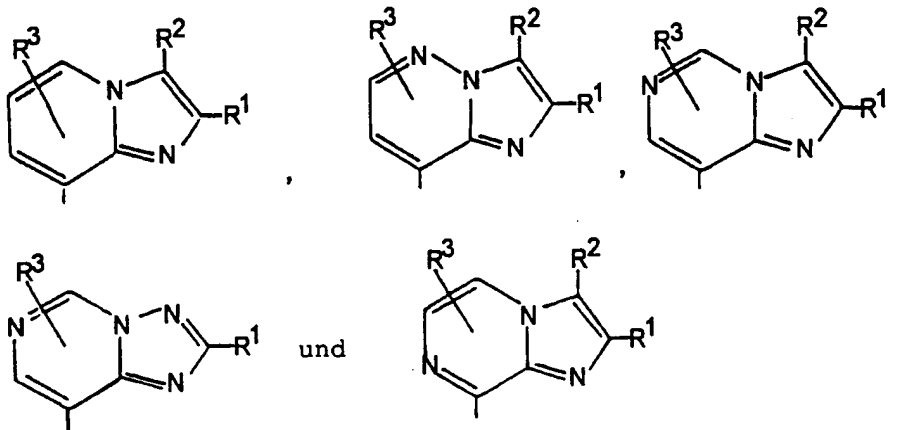
Patentansprüche für folgende Vertragsstaaten : AT, BE, CH, DE, DK, FR, GB, GR, IT, LI, LU, NL, SE

1. Kondensierte heterocyclische Verbindung der Formel:



oder ein landwirtschaftlich verträgliches Salz davon, worin:

- (1) Q eine kondensierte heterocyclische Gruppe mit einem Stickstoffatom im Brückenkopf darstellt, ausgewählt aus der Gruppe, bestehend aus

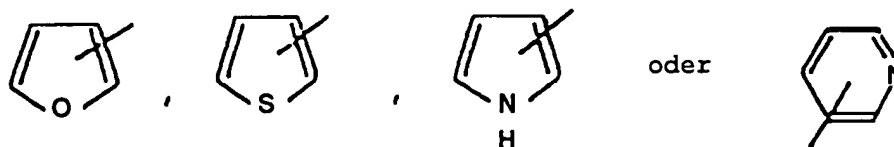


worin R<sup>1</sup> C<sub>1-6</sub>-Alkyl, Halogen, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio, C<sub>6-10</sub>-Aryloxy, C<sub>6-10</sub>-Arylthio, Alkoxycarbonyl, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein

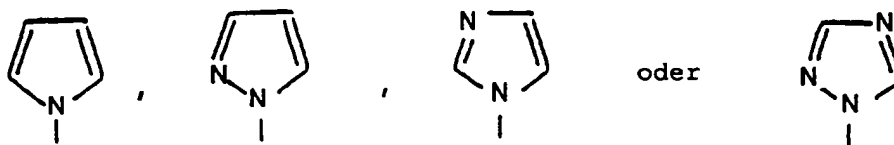
können und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder eine 5- oder 6-gliedrige, aromatische, heterocyclische Gruppe darstellt;

R<sup>2</sup> und R<sup>3</sup> Wasserstoff, C<sub>1-6</sub>-Alkyl, Halogen, Nitro, Amino, Sulfo, Mono- oder Dialkylsulfamoyl, Alkoxycarbonyl, Formyl, Cyano, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder eine 5- oder 6-gliedrige, aromatische, heterocyclische Gruppe darstellen;

X Wasserstoff; eine Gruppe, gebunden durch ein Kohlenstoffatom, ausgewählt aus der Gruppe, bestehend aus C<sub>1-10</sub>-Alkyl, C<sub>2-4</sub>-Alkenyl, C<sub>1-10</sub>-Halogenalkyl, C<sub>3-6</sub>-Cycloalkyl, C<sub>3-6</sub>-Cycloalkenyl, C<sub>6-10</sub>-Aryl, C<sub>7-10</sub>-Aralkyl und einem heterocyclischen Ring der Formel



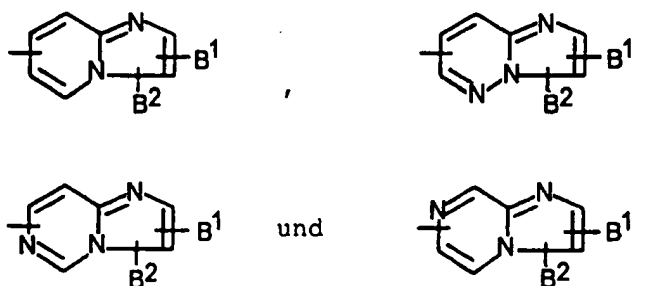
einer Gruppe, gebunden durch ein Sauerstoffatom, ausgewählt aus der Gruppe, bestehend aus C<sub>1-10</sub>-Alkoxy, C<sub>6-10</sub>-Aryloxy und C<sub>7-10</sub>-Aralkyloxy; eine Gruppe, gebunden durch ein Schwefelatom, ausgewählt aus der Gruppe, bestehend aus C<sub>1-10</sub>-Alkylthio, C<sub>6-10</sub>-Arylthio und C<sub>7-10</sub>-Aralkylthio; eine Gruppe, gebunden durch ein Stickstoffatom, ausgewählt aus der Gruppe, bestehend aus C<sub>1-10</sub>-Alkylamino, C<sub>6-10</sub>-Arylamino, C<sub>7-10</sub>-Aralkylamino und einem heterocyclischen Ring der Formel



worin jede der Gruppen, gebunden durch ein Kohlenstoffatom, ein Sauerstoffatom, ein Schwefelatom oder ein Stickstoffatom, unsubstituiert oder substituiert mit ein bis vier Substituenten ist, ausgewählt aus der Gruppe, bestehend aus Nitro, Amino, Hydroxy, Cyano, Carboxyl, Sulfonyl, Halogen, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio, Phenylthio und Benzylthio; und

Y eine Cyano-, Carbamoyl-, Thiocarbamoyl- oder Trichlormethylgruppe darstellt;

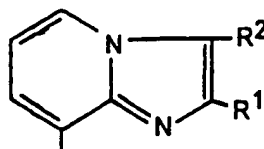
oder (2) Q eine kondensierte, heterocyclische Gruppe mit einem Stickstoffatom in dem Brückenkopf ist, ausgewählt aus der Gruppe, bestehend aus



worin die Substituenten B<sup>1</sup> und B<sup>2</sup> an der kondensierten, heterocyclischen Gruppe für Q gleich oder voneinander verschieden sein können und ein Wasserstoffatom, Nitro, Amino, Hydroxyl, Cyano, C<sub>1-3</sub>-Acyl, Carbamoyl, Carboxyl, C<sub>1-4</sub>-Alkoxycarbonyl, Sulfo, Halogen, C<sub>1-4</sub>-Alkoxy, C<sub>6-10</sub>-Aryloxy, C<sub>6-10</sub>-Arylcarbonyl, C<sub>6-10</sub>-Aryl, C<sub>7-10</sub>-Aralkyl, C<sub>3-7</sub>-Cycloalkyl, C<sub>1-4</sub>-Alkylthio, C<sub>6-10</sub>-Arylthio,

C<sub>7-10</sub>-Aralkylthio, Mono- oder Di-C<sub>1-4</sub>-alkylsulfamoyl, C<sub>1-6</sub>-Alkyl, mono- di-, tri- oder tetrahalogen-substituiertes Phenyl, mono-, di-, tri- oder tetra-C<sub>1-4</sub>-alkylsubstituiertes Phenyl, mono-, di-, tri- oder tetra-C<sub>1-4</sub>-alkoxysubstituiertes Phenyl, mono-, di-, tri- oder tetra-C<sub>1-4</sub>-alkylthiosubstituiertes Phenyl, 2-Chlor-4-nitrophenyl, 4-Nitrophenyl, 2-Methyl-4-aminophenyl, 2-Brom-4-nitrophenyl, 2-Nitro-4-methylphenyl oder eine 5- oder 6-gliedrige, heterocyclische Gruppe, ausgewählt aus Pyridyl, Furyl, Thienyl und Thiazolyl, darstellen; und X und Y die gleichen wie vorstehend definierten Bedeutungen aufweisen; oder

(3) Q ist



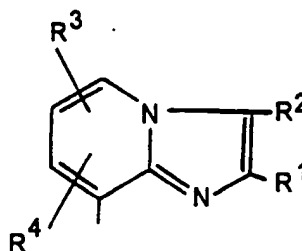
und

R<sup>1</sup> ist Methyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist 1,1,2,2-Tetrafluorethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist Morpholinomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist Isopropylthiomethyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist 2-Cyanoethyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist 2-Methyl-1-propenyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-N-Methylpyrrolyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Trifluormethylphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Methylphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 5-Methylfur-2-yl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Pyridylmethyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Methylthien-2-yl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist 1,1,2,2-Tetrafluorethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist 2-Cyanoethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist 2-Methyl-1-propenyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist 2,2,2-Trichlor-1-hydroxyethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Trifluormethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methyl-1-propenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Ethoxyiminomethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenoxyomethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenylthiomethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenylsulfinylmethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenylsulfonylmethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenoxyomethyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 3-Fluorphenyl und Y ist

Cyano;

oder R<sup>1</sup> ist Methoxymethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methoxymethyl, R<sup>2</sup> ist Morpholinomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methylthiomethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methylsulfinylmethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methylthiomethyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methylsulfonylmethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Morpholinomethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Diethoxyphosphoxymethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist (Dimethylthiocarbamoyl)thiomethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Chlormethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;

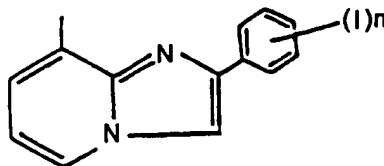
oder R<sup>1</sup> ist Trifluormethyl, R<sup>2</sup> ist Phenylcarbamoyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Wasserstoff, R<sup>2</sup> ist 1-Hydroxybenzyl, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Sulfomethyl, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 4-Dimethylaminophenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Benzylidenamino, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Biphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 4-Sulfophenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Trifluormethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Formamid;  
 oder R<sup>1</sup> ist 3-Hydroxyphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Hydroxyphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist Benzylidenamino, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Dimethylcarbamoylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methoxymethylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methyl-4-methylsulfinylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methyl-4-methylsulfonylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 4-Cyanophenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Trifluormethylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 3-Trifluormethylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methyl-5-trifluormethylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder (4) Q ist



und

R<sup>1</sup> ist Methyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, R<sup>3</sup> ist 5-Methyl, R<sup>4</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, R<sup>3</sup> ist 5-Methyl, R<sup>4</sup> ist 7-Methyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, R<sup>3</sup> ist 5-Trifluormethyl, R<sup>4</sup> ist 7-Trifluormethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, R<sup>3</sup> ist 5-Hydroxy, R<sup>4</sup> ist 6-Chlor, X ist 2-Thienyl und Y ist Cyano.

## 2. Verbindung nach Anspruch 1, worin Q

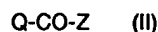


darstellt,

worin n 0, 1, 2, 3 oder 4 ist und I gleich oder verschieden ist und ausgewählt ist aus der Gruppe,  
 bestehend aus C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio oder Halogen,  
 X ist Phenyl, fluorsubstituiertes Phenyl, Thienyl oder eine Furylgruppe und  
 Y ist Cyano  
 oder deren Salz.



3. Verbindung nach Anspruch 1, nämlich  $\alpha$ -(2-Phenyl-imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)-acetonitril oder deren Salz.
4. Verbindung nach Anspruch 1, nämlich  $\alpha$ -{2-(2-Methoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitril oder deren Salz.
5. Verbindung nach Anspruch 1, nämlich  $\alpha$ -{2-(3-Methoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitril oder deren Salz.
6. Verbindung nach Anspruch 1, nämlich  $\alpha$ -{2-(2,4-Dimethoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitril oder deren Salz.
7. Verbindung nach Anspruch 1, nämlich  $\alpha$ -{2-(2,4-Diisopropylphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitril oder deren Salz.
8. Verbindung nach Anspruch 1, nämlich  $\alpha$ -{2-(3-Trifluormethylphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino}-(2-thienyl)acetonitril oder deren Salz.
9. Verfahren zur Herstellung einer kondensierten, heterocyclischen Verbindung der Formel (I) nach Anspruch 1 oder deren Salz, umfassend Umsetzen einer Verbindung der Formel

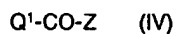


worin Q die gleiche Bedeutung wie in Anspruch 1 definiert aufweist und COZ eine Carboxylgruppe oder deren reaktives Derivat oder deren Salz bedeutet, mit einer Verbindung der allgemeinen Formel

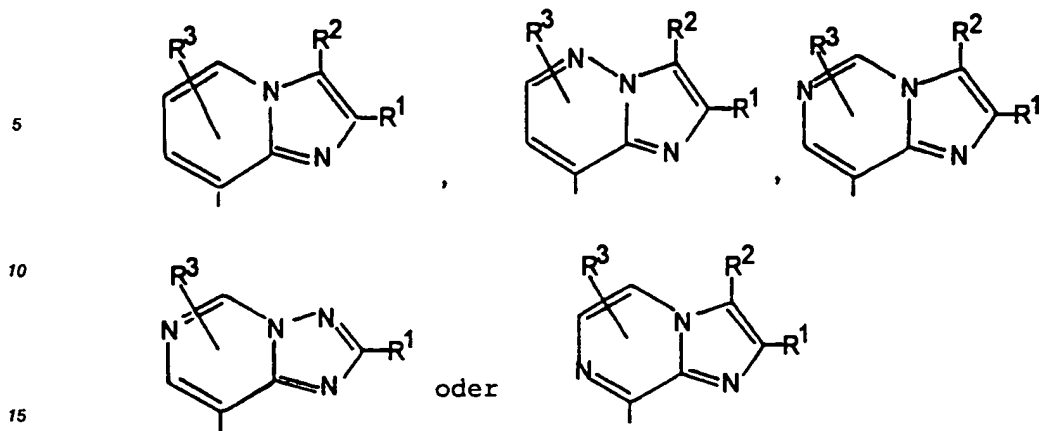


worin X und Y jeweils wie in Anspruch 1 definiert sind oder deren Salz.

10. Fungizide Zusammensetzung, umfassend eine kondensierte, heterocyclische Verbindung der Formel (I) nach Anspruch 1 oder deren Salz und ein Träger- oder Verdünnungsmittel.
11. Kondensierte, heterocyclische Carbonsäureverbindung der Formel



worin Q' eine kondensierte, heterocyclische Gruppe mit einem Stickstoffatom in dem Brückenkopf ist, wie nachstehend



20 worin R<sup>1</sup> C<sub>1-6</sub>-Alkyl, Halogen, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio, C<sub>6-10</sub>-Aryloxy, C<sub>6-10</sub>-Arylthio, Alkoxy-carbonyl, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder eine 5- oder 6-gliedrige, aromatische, heterocyclische Gruppe bedeutet;

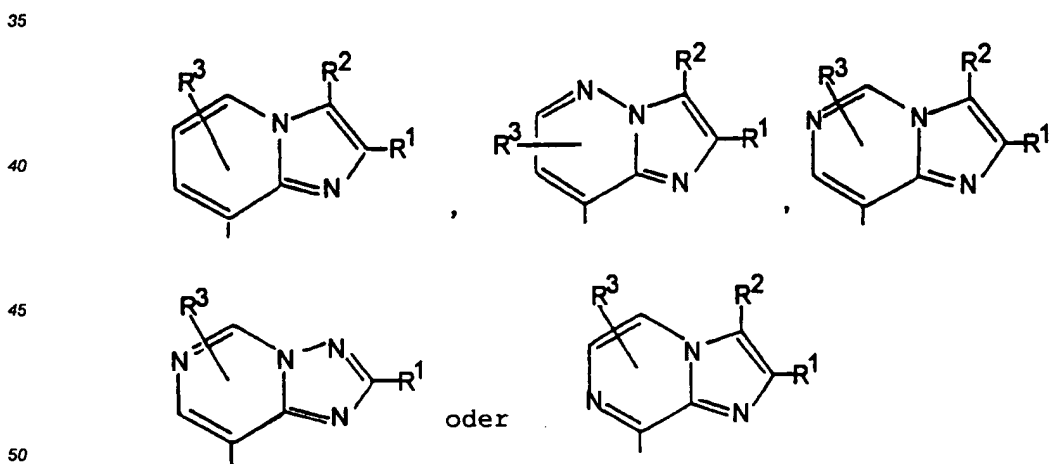
25 R<sup>2</sup> und R<sup>3</sup> Wasserstoff, C<sub>1-6</sub>-Alkyl, Halogen, Nitro, Amino, Sulfo, Mono- oder Dialkylsulfamoyl, Alkoxy-carbonyl, Formyl, Cyano, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können, und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder eine 5- oder 6-gliedrige, aromatische, heterocyclische Gruppe bedeuten

und COZ eine Carboxylgruppe oder deren reaktives Derivat bedeutet; oder deren Salz.

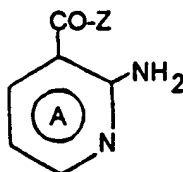
## 12. Verfahren zur Herstellung einer Verbindung der Formel

30 Q<sup>2</sup>-CO-Z (VIII)

worin Q<sup>2</sup> eine Gruppe der allgemeinen Formel



55 darstellt, worin R<sup>1</sup>, R<sup>2</sup> und R<sup>3</sup> die gleichen Bedeutungen wie in Anspruch 11 definiert aufweisen oder deren Salz, umfassend Umsetzen einer Verbindung der Formel



5

10 worin der Ring  $\textcircled{A}$  eine Pyridin-, Pyridazin-, Pyrimidin- oder Pyrazingruppe darstellt, unsubstituiert oder substituiert mit  $\text{C}_1$ - $\text{C}_6$ -Alkyl, Halogen, Nitro, Amino, Sulfo, Mono- oder Dialkylsulfamoyl, Alkoxy-carbonyl, Formyl, Cyano, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können und ausgewählt sind aus der Gruppe, bestehend aus Halogen,  $\text{C}_1$ - $\text{C}_4$ -Alkyl,  $\text{C}_1$ - $\text{C}_4$ -Alkoxy,  $\text{C}_1$ - $\text{C}_4$ -Alkylthio und Nitro oder einer 5- oder 6-gliedrigen aromatischen, heterocyclischen Gruppe und COZ eine Carboxylgruppe oder deren reaktives Derivat oder deren Salz bedeutet, mit einer Verbindung

15 der Formel:



20

oder



25

worin W ein Halogenatom darstellt und  $\text{R}^1$  und  $\text{R}^2$  die gleichen, wie vorstehend definierten Bedeutungen aufweisen.

30 13. Verwendung einer Verbindung nach den Ansprüchen 1 bis 8 als Fungizid, insbesondere zur Bekämpfung verschiedener Arten von Pflanzenkrankheiten, wie Mehltau und Brand.

#### Patentansprüche für folgenden Vertragsstaat : ES

35 1. Verfahren zur Herstellung einer kondensierten heterocyclischen Verbindung der Formel (I):



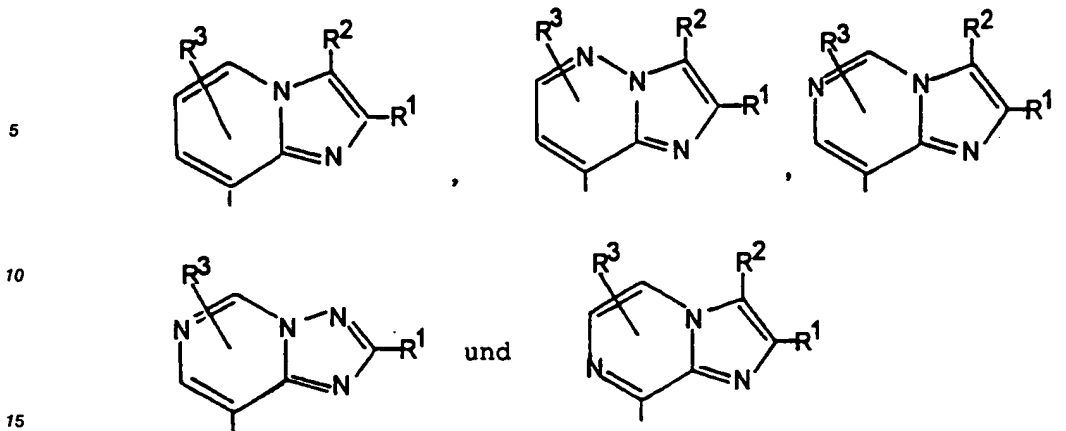
40

oder ein landwirtschaftlich verträgliches Salz davon, worin:

45 (1) Q eine kondensierte heterocyclische Gruppe mit einem Stickstoffatom im Brückenkopf darstellt, ausgewählt aus der Gruppe, bestehend aus

50

55



worin R<sup>1</sup> C<sub>1-6</sub>-Alkyl, Halogen, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio, C<sub>6-10</sub>-Aryloxy, C<sub>6-10</sub>-Arylthio, Alkoxycarbonyl, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder eine 5- oder 6-gliedrige, aromatische, heterocyclische Gruppe darstellt;

R<sup>2</sup> und R<sup>3</sup> Wasserstoff, C<sub>1-6</sub>-Alkyl, Halogen, Nitro, Amino, Sulfo, Mono- oder Dialkylsulfamoyl, Alkoxycarbonyl, Formyl, Cyano, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder eine 5- oder 6-gliedrige, aromatische, heterocyclische Gruppe darstellen;

X Wasserstoff; eine Gruppe, gebunden durch ein Kohlenstoffatom, ausgewählt aus der Gruppe, bestehend aus C<sub>1-10</sub>-Alkyl, C<sub>2-4</sub>-Alkenyl, C<sub>1-10</sub>-Halogenalkyl, C<sub>3-6</sub>-Cycloalkyl, C<sub>3-6</sub>-Cycloalkenyl, C<sub>6-10</sub>-Aryl, C<sub>7-10</sub>-Aralkyl und einem heterocyclischen Ring der Formel

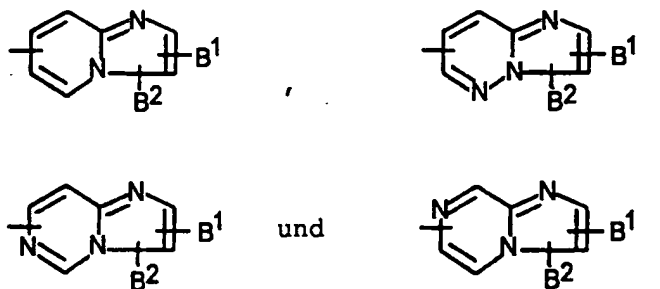


einer Gruppe, gebunden durch ein Sauerstoffatom, ausgewählt aus der Gruppe, bestehend aus C<sub>1-10</sub>-Alkoxy, C<sub>6-10</sub>-Aryloxy und C<sub>7-10</sub>-Aralkyloxy; eine Gruppe, gebunden durch ein Schwefelatom, ausgewählt aus der Gruppe, bestehend aus C<sub>1-10</sub>-Alkylthio, C<sub>6-10</sub>-Arylthio und C<sub>7-10</sub>-Aralkylthio; eine Gruppe, gebunden durch ein Stickstoffatom, ausgewählt aus der Gruppe, bestehend aus C<sub>1-10</sub>-Alkylamino, C<sub>6-10</sub>-Arylamino, C<sub>7-10</sub>-Aralkylamino und einem heterocyclischen Ring der Formel



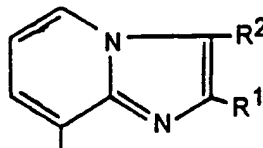
worin jede der Gruppen, gebunden durch ein Kohlenstoffatom, ein Sauerstoffatom, ein Schwefelatom oder ein Stickstoffatom, unsubstituiert oder substituiert mit ein bis vier Substituenten ist, ausgewählt aus der Gruppe, bestehend aus Nitro, Amino, Hydroxy, Cyano, Carboxyl, Sulfonyl, Halogen, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio, Phenylthio und Benzylthio; und

Y eine Cyano-, Carbamoyl-, Thiocarbamoyl- oder Trichlormethylgruppe darstellt;  
 oder (2) Q eine kondensierte, heterocyclische Gruppe mit einem Stickstoffatom in dem Brückenkopf  
 ist, ausgewählt aus der Gruppe, bestehend aus



worin die Substituenten B<sup>1</sup> und B<sup>2</sup> an der kondensierten, heterocyclischen Gruppe für Q gleich  
 oder voneinander verschieden sein können und ein Wasserstoffatom, Nitro, Amino, Hydroxyl, Cyano,  
 C<sub>1-3</sub>-Acyl, Carbamoyl, Carboxyl, C<sub>1-4</sub>-Alkoxycarbonyl, Sulfo, Halogen, C<sub>1-4</sub>-Alkoxy, C<sub>6-10</sub>-Aryloxy,  
 C<sub>6-10</sub>-Arylcarbonyl, C<sub>6-10</sub>-Aryl, C<sub>7-10</sub>-Aralkyl, C<sub>3-7</sub>-Cycloalkyl, C<sub>1-4</sub>-Alkylthio, C<sub>6-10</sub>-Arylthio,  
 C<sub>7-10</sub>-Aralkylthio, Mono- oder Di-C<sub>1-4</sub>-alkylsulfamoyl, C<sub>1-6</sub>-Alkyl, mono- di-, tri- oder tetrahalogen-  
 substituiertes Phenyl, mono-, di-, tri- oder tetra-C<sub>1-4</sub>-alkylsubstituiertes Phenyl, mono-, di-, tri- oder  
 tetra-C<sub>1-4</sub>-alkoxysubstituiertes Phenyl, mono-, di-, tri- oder tetra-C<sub>1-4</sub>-alkylthiosubstituiertes Phenyl,  
 2-Chlor-4-nitrophenyl, 4-Nitrophenyl, 2-Methyl-4-aminophenyl, 2-Brom-4-nitrophenyl, 2-Nitro-4-me-  
 thylphenyl oder eine 5- oder 6-gliedrige, heterocyclische Gruppe, ausgewählt aus Pyridyl, Furyl,  
 Thienyl und Thiazolyl, darstellen; und X und Y die gleichen wie vorstehend definierten Bedeutungen  
 aufweisen; oder

(3) Q ist

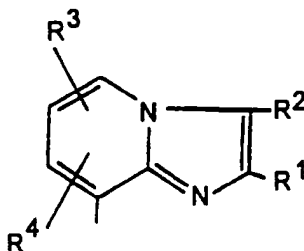


und

R<sup>1</sup> ist Methyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist 1,1,2,2-Tetrafluorethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist Morpholinomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist Isopropylthiomethyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist 2-Cyanoethyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist 2-Methyl-1-propenyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-N-Methylpyrrolyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Trifluormethylphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Methylphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 5-Methylfur-2-yl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Pyridylmethyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Methylthien-2-yl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist 1,1,2,2-Tetrafluorethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist 2-Cyanoethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist 2-Methyl-1-propenyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist 2,2,2-Trichlor-1-hydroxyethyl, X ist 2-Furyl und Y ist Cyano;

oder R<sup>1</sup> ist Trifluormethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methyl-1-propenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Ethoxyiminomethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenoxyomethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenylthiomethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenylsulfinylmethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenylsulfonylmethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenoxyomethyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 3-Fluorphenyl und Y ist Cyano;

oder R<sup>1</sup> ist Methoxymethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methoxymethyl, R<sup>2</sup> ist Morpholinomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methylthiomethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methylsulfinylmethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methylthiomethyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methylsulfonylmethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Morpholinomethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Diethoxyphosphoxymethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist (Dimethylthiocarbonyl)thiomethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Chlormethyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Furyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Trifluormethyl, R<sup>2</sup> ist Phenylcarbamoyl, X ist 3-Fluorphenyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Wasserstoff, R<sup>2</sup> ist 1-Hydroxybenzyl, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Sulfomethyl, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 4-Dimethylaminophenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Benzylidenamino, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Biphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 4-Sulfophenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Trifluormethyl, R<sup>2</sup> ist Wasserstoff, X ist 3-Fluorphenyl und Y ist Formamid;  
 oder R<sup>1</sup> ist 3-Hydroxyphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Hydroxyphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist Methyl, R<sup>2</sup> ist Benzylidenamino, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Dimethylcarbamoylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methoxymethylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methyl-4-methylsulfinylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methyl-4-methylsulfonylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 4-Cyanophenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Trifluormethylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 3-Trifluormethylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder R<sup>1</sup> ist 2-Methyl-5-trifluormethylphenyl, R<sup>2</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;  
 oder (4) Q ist



und

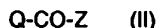
R<sup>1</sup> ist Methyl, R<sup>2</sup> ist N,N-Dimethylaminomethyl, R<sup>3</sup> ist 5-Methyl, R<sup>4</sup> ist Wasserstoff, X ist 2-Thienyl und Y ist Cyano;

oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, R<sup>3</sup> ist 5-Methyl, R<sup>4</sup> ist 7-Methyl, X ist 2-Furyl und Y ist Cyano;

oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, R<sup>3</sup> ist 5-Trifluormethyl, R<sup>4</sup> ist 7-Trifluormethyl, X ist 2-Furyl und Y ist Cyano;

oder R<sup>1</sup> ist Phenyl, R<sup>2</sup> ist Wasserstoff, R<sup>3</sup> ist 5-Hydroxy, R<sup>4</sup> ist 6-Chlor, X ist 2-Thienyl und Y ist

Cyano, umfassend Umsetzen einer Verbindung der Formel

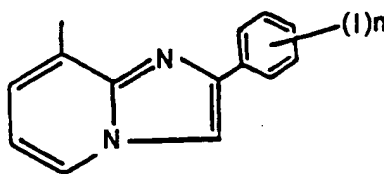


worin Q die gleiche Bedeutung wie vorstehend definiert aufweist und COZ eine Carboxylgruppe oder deren reaktives Derivat oder deren Salz bedeutet, mit einer Verbindung der allgemeinen Formel



worin X und Y jeweils wie in Anspruch 1 definiert sind oder deren Salz.

## 2. Verfahren nach Anspruch 1, worin Q



darstellt,

worin n 0, 1, 2, 3 oder 4 ist und I gleich oder verschieden ist und ausgewählt ist aus der Gruppe, bestehend aus C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio oder Halogen,

X ist Phenyl, fluorsubstituiertes Phenyl, Thienyl oder eine Furylgruppe und

Y ist Cyano oder deren Salz.

3. Verfahren zur Herstellung einer Verbindung nach Anspruch 1, nämlich  $\alpha$ -(2-Phenylimidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)acetonitril oder deren Salz.

4. Verfahren zur Herstellung einer Verbindung nach Anspruch 1, nämlich  $\alpha$ -(2-(2-Methoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)acetonitril oder deren Salz.

5. Verfahren zur Herstellung einer Verbindung nach Anspruch 1, nämlich  $\alpha$ -(2-(3-Methoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)acetonitril oder deren Salz.

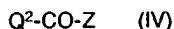
6. Verfahren zur Herstellung einer Verbindung nach Anspruch 1, nämlich  $\alpha$ -(2-(2,4-Dimethoxyphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)acetonitril oder deren Salz.

7. Verfahren zur Herstellung einer Verbindung nach Anspruch 1, nämlich  $\alpha$ -(2-(2,4-Diisopropylphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)acetonitril oder deren Salz.

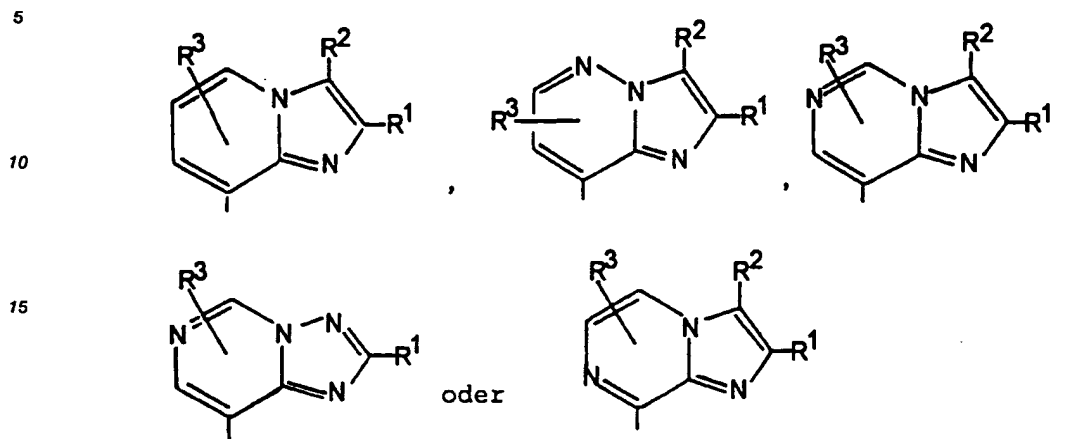
8. Verfahren zur Herstellung einer Verbindung nach Anspruch 1, nämlich  $\alpha$ -(2-(3-Trifluormethylphenyl)imidazo[1,2-a]pyridin-8-ylcarbonylamino)-(2-thienyl)acetonitril oder deren Salz.

9. Verfahren zur Herstellung einer fungiziden Zusammensetzung, umfassend Vermischen einer kondensierten heterocyclischen Verbindung der Formel (I) in Anspruch 1 oder deren Salz und ein Träger- oder Verdünnungsmittel.

10. Verfahren zur Herstellung einer Verbindung der Formel



worin Q<sup>2</sup> eine kondensierte, heterocyclische Gruppe mit einem Stickstoffatom in dem Brückenkopf ist, wie nachstehend



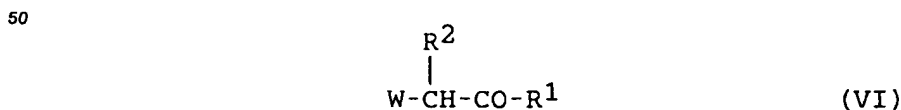
25 worin R<sup>1</sup> C<sub>1-6</sub>-Alkyl, Halogen, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio, C<sub>6-10</sub>-Aryloxy, C<sub>6-10</sub>-Arylthio, Alkoxy-carbonyl, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder eine 5- oder 6-gliedrige, aromatische, heterocyclische Gruppe bedeuten;

30 R<sup>2</sup> und R<sup>3</sup> Wasserstoff, C<sub>1-6</sub>-Alkyl, Halogen, Nitro, Amino, Sulfo, Mono- oder Dialkylsulfamoyl, Alkoxy-carbonyl, Formyl, Cyano, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können, und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder eine 5- oder 6-gliedrige, aromatische, heterocyclische Gruppe bedeuten

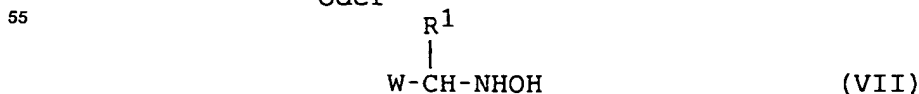
35 und COZ eine Carboxylgruppe oder deren reaktives Derivat bedeutet; oder deren Salz, umfassend Umsetzen einer Verbindung der Formel



50 worin der Ring A eine Pyridin-, Pyridazin-, Pyrimidin- oder Pyrazingruppe darstellt, unsubstituiert oder substituiert mit C<sub>1-6</sub>-Alkyl, Halogen, Nitro, Amino, Sulfo, Mono- oder Dialkylsulfamoyl, Alkoxy-carbonyl, Formyl, Cyano, Phenyl, Phenyl, substituiert mit 1 bis 4 Gruppe(n), die gleich oder verschieden sein können und ausgewählt sind aus der Gruppe, bestehend aus Halogen, C<sub>1-4</sub>-Alkyl, C<sub>1-4</sub>-Alkoxy, C<sub>1-4</sub>-Alkylthio und Nitro oder einer 5- oder 6-gliedrigen aromatischen, heterocyclischen Gruppe und COZ eine Carboxylgruppe oder deren reaktives Derivat oder deren Salz bedeutet, mit einer Verbindung der Formel:



oder





worin W ein Halogenatom darstellt und R<sup>1</sup> und R<sup>2</sup> die gleichen, wie vorstehend definierten Bedeutungen aufweisen.

11. Utilisation d'une Verbindung nach den Ansprüchen 1 bis 8 als Fungizid, insbesondere zur Bekämpfung verschiedener Arten von Pflanzenkrankheiten, wie Mehltau und Brand.

# Revendications

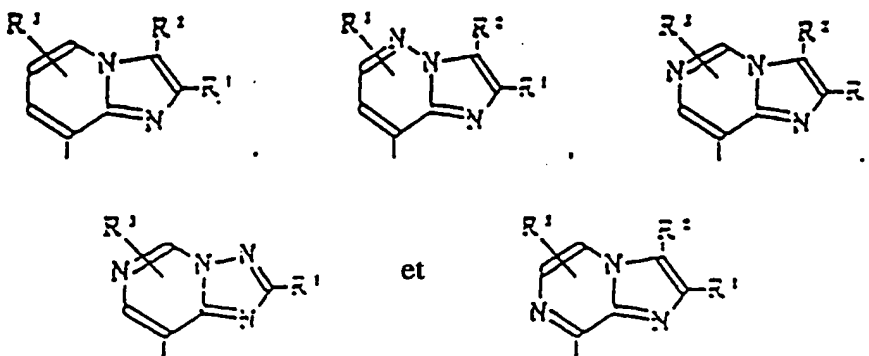
Revendications pour les Etats contractants suivants : AT, DE, CH, DE, DK, FR, GB, GR, IT, LI, LU, NL, SE

1. Composé hétérocyclique condensé, de formule :



ou sel d'un tel composé, acceptable en agriculture, où :

- 1) Q représente un groupe hétérocyclique condensé, comportant un atome d'azote en tête de pont, choisi dans l'ensemble constitué par

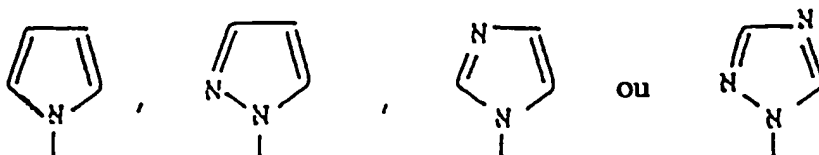


où R<sup>1</sup> représente un atome d'halogène ou un groupe alkyle en C<sub>1-6</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub>, aryloxy en C<sub>6-10</sub>, arylthio en C<sub>6-10</sub>, alcoxycarbonyl, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub> et nitro, ou encore un groupe hétérocyclique aromatique à 5 ou 6 chaînons, et R<sup>2</sup> et R<sup>3</sup> représentent des atomes d'hydrogène ou d'halogène ou des groupes alkyle en C<sub>1-6</sub>, nitro, amino, sulfo, monoalkylsulfamyle, dialkylsulfamyle, alcoxycarbonyl, formyle, cyano, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub> et nitro, ou encore des groupes hétérocycliques aromatiques à 5 ou 6 chaînons ;

X représente un atome d'hydrogène, un groupe lié par un atome de carbone et choisi dans l'ensemble constitué par les groupes alkyle en C<sub>1-10</sub>, alcényle en C<sub>2-4</sub>, halogénoalkyle en C<sub>1-10</sub>, cycloalkyle en C<sub>3-6</sub>, cycloalcényle en C<sub>3-6</sub>, aryle en C<sub>6-10</sub>, aralkyle en C<sub>7-10</sub>, et les groupes hétérocycliques de formule :



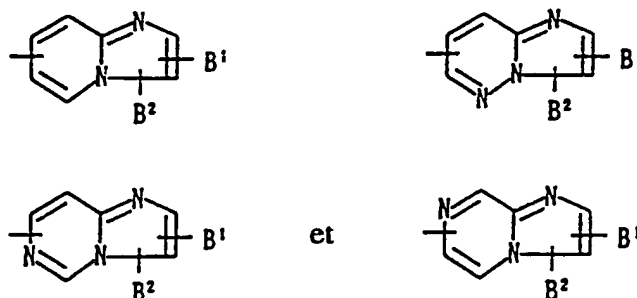
un groupe lié par un atome d'oxygène et choisi parmi les groupes alcoxy en C<sub>1-10</sub>, aryloxy en C<sub>6-10</sub> et aralcoxy en C<sub>7-10</sub>, un groupe lié par un atome de soufre et choisi parmi les groupes alkylthio en C<sub>1-10</sub>, arylthio en C<sub>6-10</sub> et aralkylthio en C<sub>7-10</sub>, ou un groupe lié par un atome d'azote et choisi parmi les groupes alkylamino en C<sub>1-10</sub>, arylamino en C<sub>6-10</sub>, aralkylamino en C<sub>7-10</sub> et les groupes hétérocycliques de formule



chacun de ces groupes liés par un atome de carbone, d'oxygène, de soufre ou d'azote portant ou non des substituants, au nombre de 1 à 4 et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes nitro, amino, hydroxy, cyano, carboxy, sulfo, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub>, phénylthio et benzylthio ; et

Y représente un groupe cyano, carbamyle, thiocarbamyle ou trichlorométhyle ; ou bien

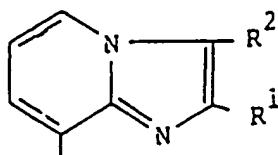
2) Q représente un groupe hétérocyclique condensé, comportant un atome d'azote en tête de pont, choisi dans l'ensemble constitué par



où les substituants B<sup>1</sup> et B<sup>2</sup> placés sur le groupe hétérocyclique condensé représenté par Q peuvent être identiques ou différents l'un de l'autre et sont des atomes d'hydrogène ou d'halogène ou des groupes nitro, amino, hydroxy, cyano, acyle en C<sub>1-3</sub>, carbamyle, carboxy, (alcoxy en C<sub>1-4</sub>)-carbonyle, sulfo, alcoxy en C<sub>1-4</sub>, aryloxy en C<sub>6-10</sub>, (aryle en C<sub>6-10</sub>)-carbonyle, aryle en C<sub>6-10</sub>, aralkyle en C<sub>7-10</sub>, cycloalkyle en C<sub>3-7</sub>, alkylthio en C<sub>1-4</sub>, arylthio en C<sub>6-10</sub>, aralkylthio en C<sub>7-10</sub>, mono- ou di-(alkyle en C<sub>1-4</sub>)-sulfamyle, alkyle en C<sub>1-5</sub>, phényle portant un, deux, trois ou quatre substituants halogéno, phényle portant un, deux, trois ou quatre substituants alkyle en C<sub>1-4</sub>, phényle portant un, deux, trois ou quatre substituants alcoxy en C<sub>1-4</sub>, phényle portant un, deux, trois ou quatre substituants alkylthio en C<sub>1-4</sub>, 2-chloro-4-nitrophényle, 4-nitrophényle, 2-méthyl-4-aminophényle, 2-bromo-4-nitrophényle ou 2-nitro-4-méthylphényle, ou encore des groupes hétérocycliques à cinq ou six chaînons, choisis parmi les groupes pyridyle, furyle, thiényle et thiazolyle ; et

X et Y ont les mêmes significations que celles qui sont indiquées plus haut ; ou bien

3) Q représente



et R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano;

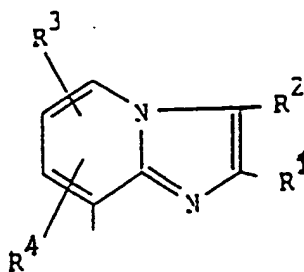
ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe 1,1,2,2-tétrafluoroéthyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
5 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe morpholinométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
10 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe isopropylthiométhyle, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe 2-cyanoéthyle, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe 2-méthyl-1-propényle, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
15 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe N-méthyl-2-pyrrolyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-trifluorométhyl-phényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 20 3-méthylphényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 5-méthyl-2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-pyridyl-méthyle et Y représente un groupe cyano ;  
25 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-méthyl-2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X 30 représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe 1,1,2,2-tétrafluoroéthyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
35 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe 2-cyanoéthyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe 2-méthyl-1-propényle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe 2,2,2-trichloro-1-hydroxyéthyle, X 40 représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe trifluorométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 2-méthyl-1-propényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
45 ou R<sup>1</sup> représente un groupe éthoxyiminométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phénoxyméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phénylthiométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
50 ou R<sup>1</sup> représente un groupe phénylsulfinylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phénylsulfonylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
55 ou R<sup>1</sup> représente un groupe phénoxyméthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthoxyméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;

ou R<sup>1</sup> représente un groupe méthoxyméthyle, R<sup>2</sup> représente un groupe morpholinométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthylthiométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
5 ou R<sup>1</sup> représente un groupe méthylsulfinylméthyle R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthylthiométhyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
10 ou R<sup>1</sup> représente un groupe méthylsulfonylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe morpholinométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe diéthoxyphosphorylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
15 ou R<sup>1</sup> représente un groupe (diméthylthiocarbamyl)thiométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe chlorométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
20 ou R<sup>1</sup> représente un groupe trifluorométhyle, R<sup>2</sup> représente un groupe phénylcarbamyle, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un atome d'hydrogène, R<sup>2</sup> représente un groupe 1-hydroxybenzyle, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe sulfométhyle, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
25 ou R<sup>1</sup> représente un groupe 4-diméthylamino-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe benzylidèneamino, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
30 ou R<sup>1</sup> représente un groupe biphenylyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 4-sulfophényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe trifluorométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
35 ou R<sup>1</sup> représente un groupe 3-hydroxyphényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 2-hydroxyphényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
40 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe benzylidèneamino, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 2-diméthylcarbamil-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 2-méthoxyméthyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
45 ou R<sup>1</sup> représente un groupe 2-méthyl-4-méthylsulfinyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 2-méthyl-4-méthylsulfonyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 4-cyanophényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
50 ou R<sup>1</sup> représente un groupe 2-trifluorométhyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 3-trifluorométhyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
55 ou R<sup>1</sup> représente un groupe 2-méthyl-5-trifluorométhyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ; ou bien

## 4) Q représente

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et R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, R<sup>3</sup> représente un groupe 5-méthyle, R<sup>4</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;

ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, R<sup>3</sup> représente un groupe 5-méthyle, R<sup>4</sup> représente un groupe 7-méthyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;

20

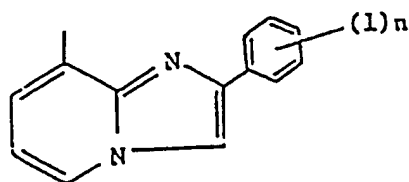
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, R<sup>3</sup> représente un groupe 5-trifluorométhyle, R<sup>4</sup> représente un groupe 7-trifluorométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;

ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, R<sup>3</sup> représente un groupe 5-hydroxy, R<sup>4</sup> représente un atome 6-chloro, X représente un groupe 2-thiényle et Y représente un groupe cyano.

25

## 2. Composé conforme à la revendication 1, dans lequel Q représente

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où n vaut 0, 1, 2, 3 ou 4, et (1) représente des substituants identiques ou différents et choisis dans l'ensemble constitué par les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub> et alkylthio en C<sub>1-4</sub>,

40

X représente un groupe phényle, phényle portant un ou des atomes de fluor en tant que substituants, thiényle ou furyle, et

Y représente un groupe cyano, ou sel d'un tel composé.

45

3. Composé conforme à la revendication 1, qui est l' $\alpha$ -(2-phényl-imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényle)acétonitrile ou l'un de ses sels.4. Composé conforme à la revendication 1, qui est l' $\alpha$ -(2-(2-méthoxyphényl)imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényle)acétonitrile ou l'un de ses sels.

50

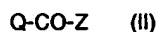
5. Composé conforme à la revendication 1, qui est l' $\alpha$ -(2-(3-méthoxyphényl)imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényle)acétonitrile ou l'un de ses sels.

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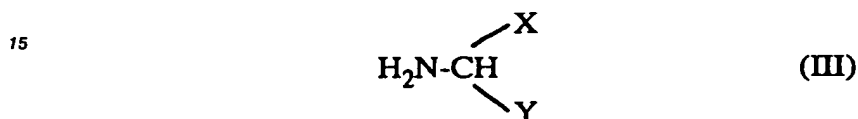
6. Composé conforme à la revendication 1, qui est l' $\alpha$ -(2-(2,4-diméthoxyphényl)imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényle)acétonitrile ou l'un de ses sels.7. Composé conforme à la revendication 1, qui est l' $\alpha$ -(2-(2,4-diisopropylphényl)imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényle)acétonitrile ou l'un de ses sels.

8. Composé conforme à la revendication 1, qui est l' $\alpha$ -{2-(3-trifluorométhylphényl)imidazo[1,2-a]pyridine-8-yl-carbonylamino}-(2-thiényle)acétonitrile ou l'un de ses sels.

9. Procédé de préparation d'un composé hétérocyclique condensé de formule (I) indiquée dans la revendication 1, ou d'un sel d'un tel composé, qui comporte le fait de faire réagir un composé de formule



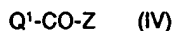
10 dans laquelle Q possède les mêmes significations que celles indiquées dans la revendication 1 et CO-Z désigne le groupe carboxy ou l'un de ses dérivés réactifs, ou un sel d'un tel composé, avec un composé de formule générale



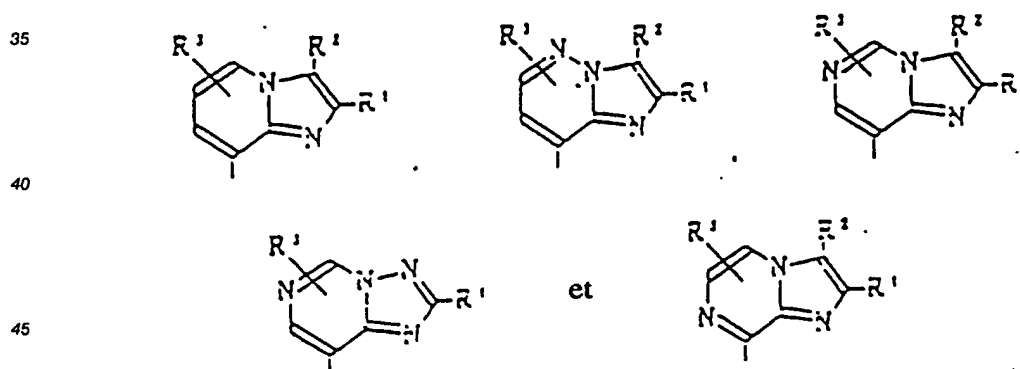
20 dans laquelle X et Y possèdent chacun les mêmes significations que celles indiquées dans la revendication 1, ou avec un sel d'un tel composé.

10. Composition fongicide comprenant un composé hétérocyclique condensé de formule (I), conforme à la revendication 1, ou un sel d'un tel composé, et un véhicule ou un diluant.

11. Composé hétérocyclique condensé à fonction de type acide carboxylique, de formule



30 dans laquelle Q<sup>1</sup> représente un groupe hétérocyclique condensé, comportant un atome d'azote en tête de pont, choisi dans l'ensemble constitué par



50 où R<sup>1</sup> représente un atome d'halogène ou un groupe alkyle en C<sub>1-6</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub>, aryloxy en C<sub>6-10</sub>, arylthio en C<sub>6-10</sub>, alcoxycarbonyl, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub> et nitro, ou encore un groupe hétérocyclique aromatique à 5 ou 6 chaînons, et

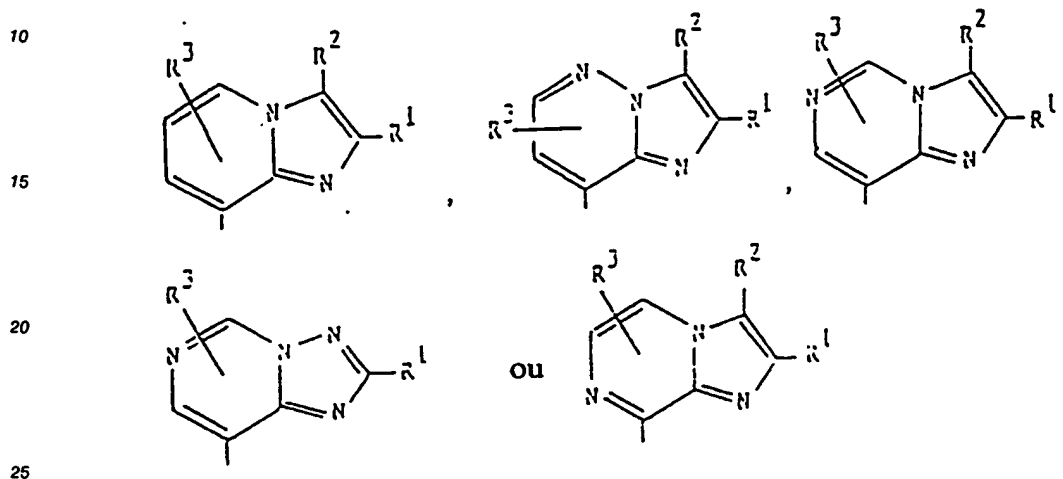
55 R<sup>2</sup> et R<sup>3</sup> représentent des atomes d'hydrogène ou d'halogène ou des groupes alkyle en C<sub>1-6</sub>, nitro, amino, sulfo, monoalkylsulfamyle, dialkylsulfamyle, alcoxycarbonyl, formyle, cyano, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub> et nitro, ou encore des groupes hétérocycliques aromatiques à 5 ou 6 chaînons,

et -COZ désigne le groupe carboxy ou l'un de ses dérivés réactifs, ou sel d'un tel composé.

## 12. Procédé de préparation d'un composé de formule

5 Q<sup>2</sup>-CO-Z (VIII)

dans laquelle Q<sup>2</sup> représente un groupe de formule générale

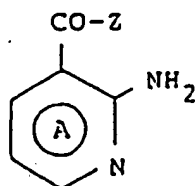


dans laquelle R<sup>1</sup>, R<sup>2</sup> et R<sup>3</sup> ont les mêmes significations que celles qui sont indiquées dans la revendication 11,

ou d'un sel d'un tel composé, lequel procédé comporte le fait de faire réagir un composé de formule

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dans laquelle le cycle A est un cycle de type pyridine, pyridazine, pyrimidine ou pyrazine, qui porte ou non, en tant que substituants, des atomes d'halogène ou des groupes alkyle en C<sub>1-5</sub>, nitro, amino, sulfo, monoalkylsulfamyle, dialkylsulfamyle, alcoxycarbonyle, formyle, cyano, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub> et nitro, ou encore

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des groupes hétérocycliques aromatiques à 5 ou 6 chaînons, et -COZ désigne le groupe carboxy ou l'un de ses dérivés réactifs,

ou un sel d'un tel composé, avec un composé de formule

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ou

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où W représente un atome d'halogène et R<sup>1</sup> et R<sup>2</sup> ont les mêmes significations que celles qui sont indiquées plus haut.

13. Emploi d'un composé conforme à l'une des revendications 1 à 8 en tant qu'agent fongicide, et plus particulièrement, dans la lutte contre divers types de flétrissure généralisée des plantes, comme le mildiou et la rouille des céréales.

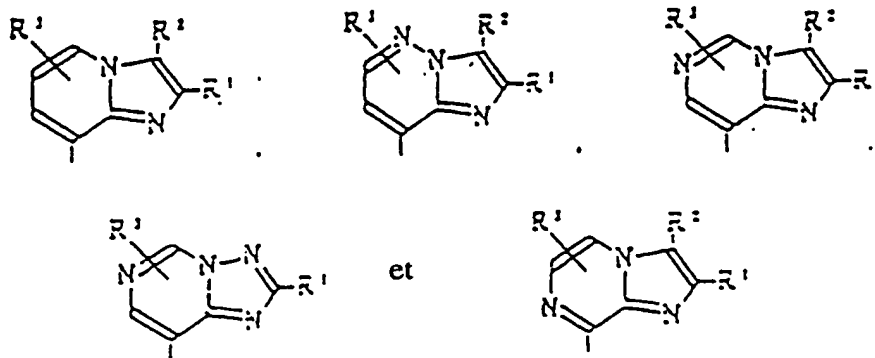
#### Revendications pour l'Etat contractant suivant : ES

1. Procédé de préparation d'un composé hétérocyclique condensé, de formule :



ou d'un sel d'un tel composé, acceptable en agriculture, où :

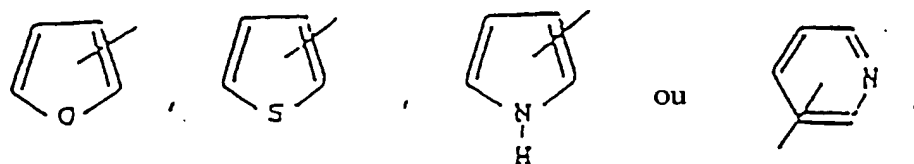
- 1) Q représente un groupe hétérocyclique condensé, comportant un atome d'azote en tête de pont, choisi dans l'ensemble constitué par



où R<sup>1</sup> représente un atome d'halogène ou un groupe alkyle en C<sub>1-6</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub>, aryloxy en C<sub>6-10</sub>, arylthio en C<sub>6-10</sub>, alcoxycarbonyl, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub> et nitro, ou encore un groupe hétérocyclique aromatique à 5 ou 6 chaînons, et

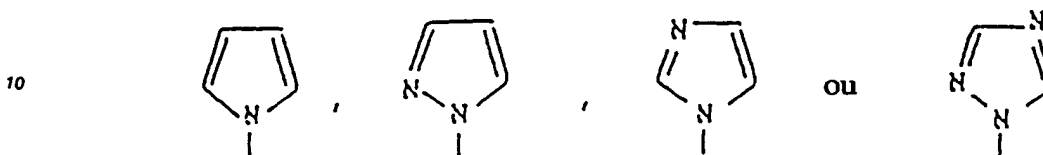
R<sup>2</sup> et R<sup>3</sup> représentent des atomes d'hydrogène ou d'halogène ou des groupes alkyle en C<sub>1-6</sub>, nitro, amino, sulfo, monoalkylsulfamyle, dialkylsulfamyle, alcoxycarbonyl, formyle, cyano, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub> et nitro, ou encore des groupes hétérocycliques aromatiques à 5 ou 6 chaînons ;

X représente un atome d'hydrogène, un groupe lié par un atome de carbone et choisi dans l'ensemble constitué par les groupes alkyle en C<sub>1-10</sub>, alcényle en C<sub>2-4</sub>, halogénoalkyle en C<sub>1-10</sub>, cycloalkyle en C<sub>3-6</sub>, cycloalcényle en C<sub>3-6</sub>, aryle en C<sub>6-10</sub>, aralkyle en C<sub>7-10</sub>, et les groupes hétérocycliques de formule :





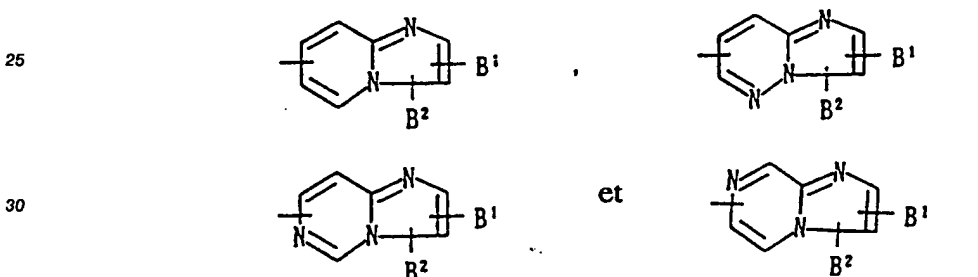
un groupe lié par un atome d'oxygène et choisi parmi les groupes alcoxy en C<sub>1-10</sub>, aryloxy en C<sub>6-10</sub> et aralcoxy en C<sub>7-10</sub>, un groupe lié par un atome de soufre et choisi parmi les groupes alkylthio en C<sub>1-10</sub>, arylthio en C<sub>6-10</sub> et aralkylthio en C<sub>7-10</sub>, ou un groupe lié par un atome d'azote et choisi parmi les groupes alkylamino en C<sub>1-10</sub>, arylamino en C<sub>6-10</sub>, aralkylamino en C<sub>7-10</sub> et les groupes hétérocycliques de formule



15 chacun de ces groupes liés par un atome de carbone, d'oxygène, de soufre ou d'azote portant ou non des substituants, au nombre de 1 à 4 et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes nitro, amino, hydroxy, cyano, carboxy, sulfo, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub>, phénylthio et benzylthio ; et

Y représente un groupe cyano, carbamyle, thiocarbamyle ou trichlorométhyle ; ou bien

2) Q représente un groupe hétérocyclique condensé, comportant un atome d'azote en tête de pont, choisi dans l'ensemble constitué par



35 où les substituants B<sup>1</sup> et B<sup>2</sup> placés sur le groupe hétérocyclique condensé représenté par Q peuvent être identiques ou différents l'un de l'autre et sont des atomes d'hydrogène ou d'halogène ou des groupes nitro, amino, hydroxy, cyano, acyle en C<sub>1-3</sub>, carbamyle, carboxy, (alcoxy en C<sub>1-4</sub>)-carbonyle, sulfo, alcoxy en C<sub>1-4</sub>, aryloxy en C<sub>6-10</sub>, (aryle en C<sub>6-10</sub>)-carbonyle, aryle en C<sub>6-10</sub>, aralkyle en C<sub>7-10</sub>, cycloalkyle en C<sub>3-7</sub>, alkylthio en C<sub>1-4</sub>, arylthio en C<sub>6-10</sub>, aralkylthio en C<sub>7-10</sub>, mono- ou di-(alkyle en C<sub>1-4</sub>)-sulfamyle, alkyle en C<sub>1-6</sub>, phényle portant un, deux, trois ou quatre substituants halogéno, phényle portant un, deux, trois ou quatre substituants alkyle en C<sub>1-4</sub>, phényle portant un, deux, trois ou quatre substituants alcoxy en C<sub>1-4</sub>, phényle portant un, deux, trois ou quatre substituants alkylthio en C<sub>1-4</sub>, 2-chloro-4-nitrophényle, 4-nitrophényle, 2-méthyl-4-aminophényle, 2-bromo-4-nitrophényle ou 2-nitro-4-méthylphényle, ou encore des groupes hétérocycliques à cinq ou six chaînons, choisis parmi les groupes pyridyle, furyle, thiényle et thiazolyle ; et

X et Y ont les mêmes significations que celles qui sont indiquées plus haut ; ou bien

3) Q représente



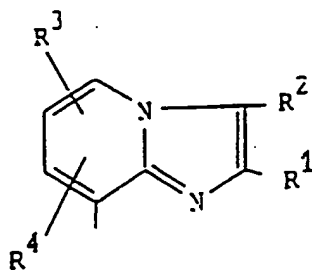
55 et R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;

ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe 1,1,2,2-tétrafluoroéthyle, X représente

un groupe 2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X  
 représente un groupe 2-thiényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X  
 5 représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe morpholinométhyle, X représente un  
 groupe 2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe isopropylthiométhyle, X représente  
 un groupe 3-fluorophényle et Y représente un groupe cyano ;  
 10 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe 2-cyanoéthyle, X représente un  
 groupe 3-fluorophényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe 2-méthyl-1-propényle, X représente  
 un groupe 3-fluorophényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe  
 15 N-méthyl-2-pyrrolyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe  
 3-trifluorométhyl-phényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe  
 3-méthylphényle et Y représente un groupe cyano ;  
 20 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe  
 5-méthyl-2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe  
 2-pyridyl-méthyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe  
 25 3-méthyl-2-thiényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X  
 représente un groupe 2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X  
 représente un groupe 2-thiényle et Y représente un groupe cyano ;  
 30 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X  
 représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe 1,1,2,2-tétrafluoroéthyle, X représente  
 un groupe 2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe 2-cyano-éthyle, X représente un  
 35 groupe 2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe 2-méthyl-1-propényle, X représente  
 un groupe 2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe 2,2,2-trichloro-1-hydroxyéthyle, X  
 représente un groupe 2-furyle et Y représente un groupe cyano ;  
 40 ou R<sup>1</sup> représente un groupe trifluorométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un  
 groupe 2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe 2-méthyl-1-propényle, R<sup>2</sup> représente un atome d'hydrogène, X repré-  
 sente un groupe 2-thiényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe éthoxyiminométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente  
 45 un groupe 2-thiényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phénoxyméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un  
 groupe 3-fluorophényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phénylthiométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente  
 un groupe 3-fluorophényle et Y représente un groupe cyano ;  
 50 ou R<sup>1</sup> représente un groupe phénylsulfinylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X repré-  
 sente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phénylsulfonylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X  
 représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe phénoxyméthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle,  
 55 X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe méthoxyméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un  
 groupe 2-furyle et Y représente un groupe cyano ;  
 ou R<sup>1</sup> représente un groupe méthoxyméthyle, R<sup>2</sup> représente un groupe morpholinométhyle, X

représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthylthiométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthylsulfinylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
5 ou R<sup>1</sup> représente un groupe méthylthiométhyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthylsulfonylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
10 ou R<sup>1</sup> représente un groupe morpholinométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe diéthoxyphosphorylméthyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe (diméthylthiocarbamyl)thiométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
15 ou R<sup>1</sup> représente un groupe chlorométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-furyle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe trifluorométhyle, R<sup>2</sup> représente un groupe phénylcarbamyle, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
20 ou R<sup>1</sup> représente un atome d'hydrogène, R<sup>2</sup> représente un groupe 1-hydroxybenzyle, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe sulfométhyle, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 4-diméthylamino-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
25 ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un groupe benzylidèneamino, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe biphénylyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
30 ou R<sup>1</sup> représente un groupe 4-sulfophényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe trifluorométhyle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 3-fluorophényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 3-hydroxyphényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
35 ou R<sup>1</sup> représente un groupe 2-hydroxyphényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe benzylidèneamino, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
40 ou R<sup>1</sup> représente un groupe 2-diméthylcarbamyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 2-méthoxyméthyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 2-méthyl-4-méthylsulfinyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
45 ou R<sup>1</sup> représente un groupe 2-méthyl-4-méthylsulfonyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 4-cyanophényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
50 ou R<sup>1</sup> représente un groupe 2-trifluorométhyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 3-trifluorométhyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;  
ou R<sup>1</sup> représente un groupe 2-méthyl-5-trifluorométhyl-phényle, R<sup>2</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ; ou bien

4) Q représente



et R<sup>1</sup> représente un groupe méthyle, R<sup>2</sup> représente un groupe N,N-diméthylaminométhyle, R<sup>3</sup> représente un groupe 5-méthyle, R<sup>4</sup> représente un atome d'hydrogène, X représente un groupe 2-thiényle et Y représente un groupe cyano ;

ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, R<sup>3</sup> représente un groupe 5-méthyle, R<sup>4</sup> représente un groupe 7-méthyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;

ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, R<sup>3</sup> représente un groupe 5-trifluorométhyle, R<sup>4</sup> représente un groupe 7-trifluorométhyle, X représente un groupe 2-furyle et Y représente un groupe cyano ;

ou R<sup>1</sup> représente un groupe phényle, R<sup>2</sup> représente un atome d'hydrogène, R<sup>3</sup> représente un groupe 5-hydroxy, R<sup>4</sup> représente un atome 6-chloro, X représente un groupe 2-thiényle et Y représente un groupe cyano, lequel procédé comporte le fait de faire réagir un composé de formule

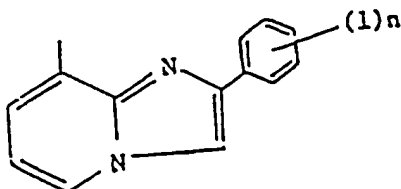
Q-CO-Z (II)

dans laquelle Q possède les mêmes significations que celles qui sont indiquées plus haut et CO-Z désigne le groupe carboxy ou l'un de ses dérivés réactifs, ou un sel d'un tel composé, avec un composé de formule générale



dans laquelle X et Y possèdent chacun les mêmes significations que celles qui sont indiquées plus haut, ou avec un sel d'un tel composé.

2. Procédé, conforme à la revendication 1, de préparation d'un composé dans lequel Q représente

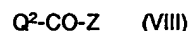


où n vaut 0, 1, 2, 3 ou 4, et (I) représente des substituants identiques ou différents et choisis dans l'ensemble constitué par les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub> et alkylthio en C<sub>1-4</sub>,

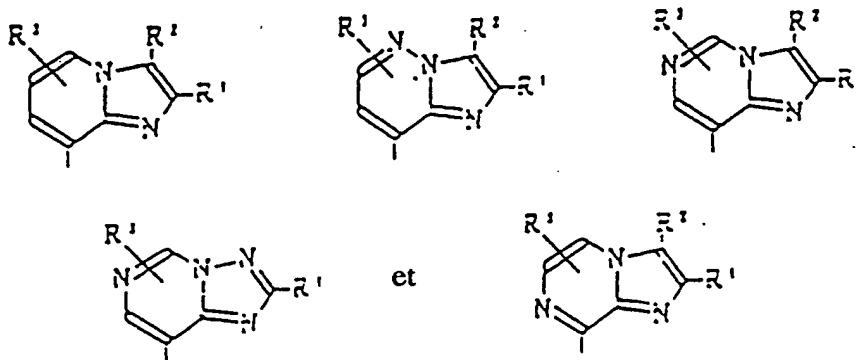
X représente un groupe phényle, phényle portant un ou des atomes de fluor en tant que substituants, thiényle ou furyle, et

Y représente un groupe cyano,  
ou d'un sel d'un tel composé.

3. Procédé, conforme à la revendication 1, de préparation d'un composé qui est l' $\alpha$ -(2-phényl-imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényl)acétonitrile ou l'un de ses sels.
4. Procédé, conforme à la revendication 1, de préparation d'un composé qui est l' $\alpha$ -(2-(2-méthoxyphényl)-imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényl)acétonitrile ou l'un de ses sels.
5. Procédé, conforme à la revendication 1, de préparation d'un composé qui est l' $\alpha$ -(2-(3-méthoxyphényl)-imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényl)acétonitrile ou l'un de ses sels.
6. Procédé, conforme à la revendication 1, de préparation d'un composé qui est l' $\alpha$ -(2-(2,4-diméthoxyphényl)-imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényl)acétonitrile ou l'un de ses sels.
7. Procédé, conforme à la revendication 1, de préparation d'un composé qui est l' $\alpha$ -(2-(2,4-diisopropylphényl)-imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényl)acétonitrile ou l'un de ses sels.
8. Procédé, conforme à la revendication 1, de préparation d'un composé qui est l' $\alpha$ -(2-(3-trifluorométhylphényl)-imidazo[1,2-a]pyridine-8-yl-carbonylamino)-(2-thiényl)acétonitrile ou l'un de ses sels.
9. Procédé de préparation d'une composition fongicide, qui comporte le fait de mélanger un composé hétérocyclique condensé de formule (I), indiquée dans la revendication 1, ou un sel d'un tel composé, et un véhicule ou un diluant.
10. Procédé de préparation d'un composé de formule



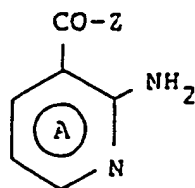
dans laquelle  $Q^2$  représente un groupe hétérocyclique condensé, comportant un atome d'azote en tête de pont, choisi dans l'ensemble constitué par



où  $R^1$  représente un atome d'halogène ou un groupe alkyle en  $C_{1-6}$ , alcoxy en  $C_{1-4}$ , alkylthio en  $C_{1-4}$ , aryloxy en  $C_{6-10}$ , arylthio en  $C_{6-10}$ , alcoxycarbonyl, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en  $C_{1-4}$ , alcoxy en  $C_{1-4}$ , alkylthio en  $C_{1-4}$  et nitro, ou encore un groupe hétérocyclique aromatique à 5 ou 6 chaînons, et

$R^2$  et  $R^3$  représentent des atomes d'hydrogène ou d'halogène ou des groupes alkyle en  $C_{1-6}$ , nitro, amino, sulfo, monoalkylsulfamyle, dialkylsulfamyle, alcoxycarbonyl, formyle, cyano, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en  $C_{1-4}$ , alcoxy en  $C_{1-4}$ , alkylthio en  $C_{1-4}$  et nitro, ou encore des groupes hétérocycliques aromatiques à 5 ou 6 chaînons ;  
et -COZ désigne le groupe carboxy ou l'un de ses dérivés réactifs ;

ou d'un sel d'un tel composé,  
lequel procédé comporte le fait de faire réagir un composé de formule

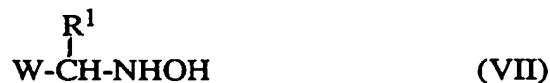


dans laquelle le cycle A est un cycle de type pyridine, pyridazine, pyrimidine ou pyrazine, qui porte ou non, en tant que substituants, des atomes d'halogène ou des groupes alkyle en C<sub>1-6</sub>, nitro, amino, sulfo, monoalkylsulfamyle, dialkylsulfamyle, alcoxycarbonyle, formyle, cyano, phényle, ou phényle portant de 1 à 4 substituants, identiques ou différents et choisis dans l'ensemble que constituent les atomes d'halogène et les groupes alkyle en C<sub>1-4</sub>, alcoxy en C<sub>1-4</sub>, alkylthio en C<sub>1-4</sub> et nitro, ou encore des groupes hétérocycliques aromatiques à 5 ou 6 chaînons, et -COZ désigne le groupe carboxy ou l'un de ses dérivés réactifs,

ou un sel d'un tel composé, avec un composé de formule



ou



où W représente un atome d'halogène et R<sup>1</sup> et R<sup>2</sup> ont les mêmes significations que celles qui sont indiquées plus haut.

11. Emploi d'un composé préparé conformément à l'une des revendications 1 à 8 en tant qu'agent fongicide, et plus particulièrement, dans la lutte contre divers types de flétrissure généralisée des plantes, comme le mildiou et la rouille des céréales.